

(12) INTERNATIONAL APPLICATION PUBLISHED UNDER THE PATENT COOPERATION TREATY (PCT)

(19) World Intellectual Property Organization
International Bureau



(43) International Publication Date
13 November 2003 (13.11.2003)

PCT

(10) International Publication Number
WO 03/093267 A1

(51) International Patent Classification⁷: **C07D 471/08, A61K 31/4995**

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(21) International Application Number: **PCT/EP03/03721**

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(22) International Filing Date: 8 April 2003 (08.04.2003)

(81) Designated States (*national*): AE, AG, AL, AM, AT, AU, AZ, BA, BB, BG, BR, BY, BZ, CA, CH, CN, CO, CR, CU, CZ, DE, DK, DM, DZ, EC, EE, ES, FI, GB, GD, GE, GH, GM, HR, HU, ID, IL, IN, IS, JP, KE, KG, KP, KR, KZ, LC, LK, LR, LS, LT, LU, LV, MA, MD, MG, MK, MN, MW, MX, MZ, NO, NZ, OM, PH, PL, PT, RO, RU, SC, SD, SE, SG, SK, SL, TJ, TM, TN, TR, TT, TZ, UA, UG, US, UZ, VC, VN, YU, ZA, ZM, ZW.

(25) Filing Language: English

(26) Publication Language: English

(30) Priority Data:
PCT/EP02/04705 29 April 2002 (29.04.2002) EP

(84) Designated States (*regional*): ARIPO patent (GH, GM, KE, LS, MW, MZ, SD, SL, SZ, TZ, UG, ZM, ZW), Eurasian patent (AM, AZ, BY, KG, KZ, MD, RU, TJ, TM), European patent (AT, BE, BG, CH, CY, CZ, DE, DK, EE, ES, FI, FR, GB, GR, HU, IE, IT, LU, MC, NL, PT, RO, SE, SI, SK, TR), OAPI patent (BF, BJ, CF, CG, CI, CM, GA, GN, GQ, GW, ML, MR, NE, SN, TD, TG).

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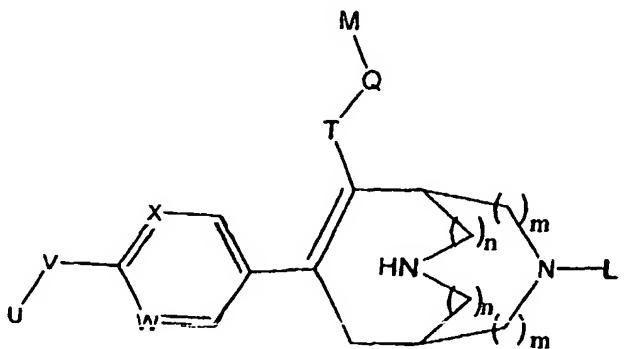
Published:

— with international search report

For two-letter codes and other abbreviations, refer to the "Guidance Notes on Codes and Abbreviations" appearing at the beginning of each regular issue of the PCT Gazette.

(54) Title: 7-ARYL-3,9-DIAZABICYCLO(3.3.1)NON-6-ENE DERIVATIVES AND THEIR USE AS RENIN INHIBITORS IN THE TREATMENT OF HYPERTENSION, CARDIOVASCULAR OR RENAL DISEASES

WO 03/093267 A1



(57) Abstract: The invention relates to novel 3,9-diazabicyclo[3.3.1]nonene derivatives of formula (I) and related compounds and their use as active ingredients in the preparation of pharmaceutical compositions. The invention also concerns related aspects including processes for the preparation of the compounds, pharmaceutical compositions containing one or more of those compounds and especially their use as inhibitors of renin.

Novel Diazabicyclononene Derivatives

5 The invention relates to novel compounds of the general formula I. The invention also concerns related aspects including processes for the preparation of the compounds, pharmaceutical compositions containing one or more compounds of formula I and especially their use as renin inhibitors in cardiovascular events and renal insufficiency. Furthermore, these compounds can be regarded as inhibitors
10 of other aspartyl proteases and might therefore be useful as inhibitors of plasmepsins to treat malaria and as inhibitors of *Candida albicans* secreted aspartyl proteases to treat fungal infections.

In the renin-angiotensin system (RAS) the biologically active angiotensin II (Ang
15 II) is generated by a two-step mechanism. The highly specific enzyme renin cleaves angiotensinogen to angiotensin I (Ang I), which is then further processed to Ang II by the less specific angiotensin-converting enzyme (ACE). Ang II is known to work on at least two receptor subtypes called AT₁ and AT₂. Whereas AT₁ seems to transmit most of the known functions of Ang II, the role of AT₂ is
20 still unknown.

Modulation of the RAS represents a major advance in the treatment of cardiovascular diseases. ACE inhibitors and AT₁ blockers have been accepted to treat hypertension (Waeber B. *et al.*, "The renin-angiotensin system: role in
25 experimental and human hypertension", in Berkenhager W. H., Reid J. L. (eds): *Hypertension*, Amsterdam, Elsevier Science Publishing Co, 1996, 489-519; Weber M. A., *Am. J. Hypertens.*, 1992, 5, 247S). In addition, ACE inhibitors are used for renal protection (Rosenberg M. E. *et al.*, *Kidney International*, 1994, 45, 403; Breyer J. A. *et al.*, *Kidney International*, 1994, 45, S156), in the prevention
30 of congestive heart failure (Vaughan D. E. *et al.*, *Cardiovasc. Res.*, 1994, 28, 159;

Fouad-Tarazi F. *et al.*, *Am. J. Med.*, 1988, 84 (*Suppl. 3A*), 83) and myocardial infarction (Pfeffer M. A. *et al.*, *N. Engl. J. Med.*, 1992, 327, 669).

The rationale to develop renin inhibitors is the specificity of renin (Kleinert H. D., 5 *Cardiovasc. Drugs*, 1995, 9, 645). The only substrate known for renin is angiotensinogen, which can only be processed (under physiological conditions) by renin. In contrast, ACE can also cleave bradykinin besides Ang I and can be bypassed by chymase, a serine protease (Husain A., *J. Hypertens.*, 1993, 11, 1155). In patients inhibition of ACE thus leads to bradykinin accumulation causing 10 cough (5-20%) and potentially life-threatening angioneurotic edema (0.1-0.2%) (Israili Z. H. *et al.*, *Annals of Internal Medicine*, 1992, 117, 234). Chymase is not inhibited by ACE inhibitors. Therefore, the formation of Ang II is still possible in patients treated with ACE inhibitors. Blockade of the AT₁ receptor (e.g. by losartan) on the other hand overexposes other AT-receptor subtypes to Ang II, 15 whose concentration is dramatically increased by the blockade of AT₁ receptors. This may raise serious questions regarding the safety and efficacy profile of AT₁ receptor antagonists. In summary, renin inhibitors are not only expected to be different from ACE inhibitors and AT₁ blockers with regard to safety, but more importantly also with regard to their efficacy to block the RAS.

20 Only limited clinical experience (Azizi M. *et al.*, *J. Hypertens.*, 1994, 12, 419; Neutel J. M. *et al.*, *Am. Heart*, 1991, 122, 1094) has been created with renin inhibitors because of their insufficient oral activity due to their peptidomimetic character (Kleinert H. D., *Cardiovasc. Drugs*, 1995, 9, 645). The clinical 25 development of several compounds has been stopped because of this problem together with the high cost of goods. Only one compound containing four chiral centers has entered clinical trials (Rahuel J. *et al.*, *Chem. Biol.*, 2000, 7, 493; Mealy N. E., *Drugs of the Future*, 2001, 26, 1139). Thus, metabolically stable, orally bioavailable and sufficiently soluble renin inhibitors that can be prepared on 30 a large scale are missing and sought. Recently, the first non-peptide renin inhibitors were described which show high *in vitro* activity (Oefner C. *et al.*, *Chem. Biol.*, 1999, 6, 127; Patent Application WO97/09311; Märki H. P. *et al.*, *II*

Farmaco, 2001, 56, 21). However, the development status of these compounds is not known.

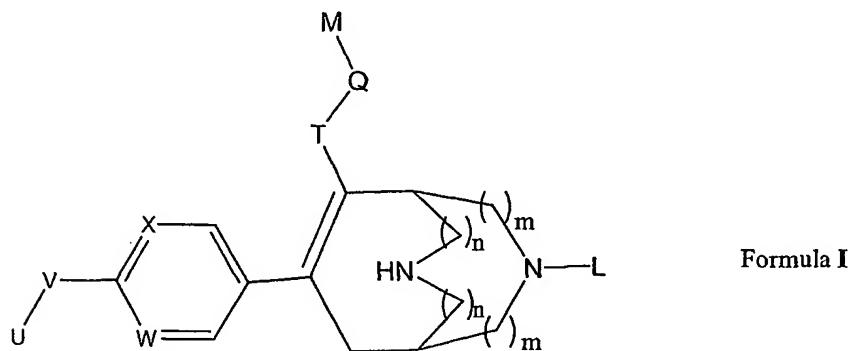
The present invention relates to the identification of renin inhibitors of a non-peptidic nature and of low molecular weight. Orally active renin inhibitors of long duration of action which are active in indications beyond blood pressure regulation where the tissular renin-chymase system may be activated leading to pathophysiologically altered local functions such as renal, cardiac and vascular remodeling, atherosclerosis, and possibly restenosis are described.

10

The present invention describes non-peptidic renin inhibitors.

In particular, the present invention relates to novel compounds of the general formula I,

15



wherein

X and W represent independently a nitrogen atom or a -CH- group;

20

V represents -(CH₂)_r; -A-(CH₂)_s; -CH₂-A-(CH₂)_t; -(CH₂)_s-A-; -(CH₂)₂-A-(CH₂)_u-; -A-(CH₂)_v-B-; -CH₂-CH₂-CH₂-A-CH₂-; -A-CH₂-CH₂-B-CH₂-; -CH₂-A-CH₂-CH₂-B-; -CH₂-CH₂-A-CH₂-CH₂-; -CH₂-CH₂-CH₂-CH₂-A-CH₂-; -A-CH₂-CH₂-B-CH₂-CH₂-; -CH₂-A-CH₂-CH₂-B-CH₂-; -CH₂-A-CH₂-CH₂-CH₂-B-; or
25 -CH₂-CH₂-A-CH₂-CH₂-B-;

A and B independently represent -O-; -S-; -SO-; -SO₂-;

U represents aryl; heteroaryl;

- 5 T represents -CONR¹-; -(CH₂)_pOCO-; -(CH₂)_pN(R¹)CO-; -(CH₂)_pN(R¹)SO₂-; or
-COO-;

Q represents lower alkylene; lower alkenylene;

- 10 M represents hydrogen; cycloalkyl; aryl; heterocyclyl; heteroaryl;

L represents -R³; -COR³; -COOR³; -CONR²R³; -SO₂R³; -SO₂NR²R³;
-COCH(Aryl)₂;

- 15 R¹ represents hydrogen; lower alkyl; lower alkenyl; lower alkinyl; cycloalkyl;
aryl; cycloalkyl - lower alkyl;

R² and R^{2'} independently represent hydrogen; lower alkyl; lower alkenyl;
cycloalkyl; cycloalkyl - lower alkyl;

20

- R³ represents hydrogen; lower alkyl; lower alkenyl; cycloalkyl; aryl; heteroaryl;
heterocyclyl; cycloalkyl - lower alkyl; aryl - lower alkyl; heteroaryl - lower alkyl;
heterocyclyl - lower alkyl; aryloxy - lower alkyl; heteroaryloxy - lower alkyl,
whereby these groups may be unsubstituted or mono-, di- or trisubstituted with
25 hydroxy, -OCOR², -COOR², lower alkoxy, cyano, -CONR²R^{2'}, CO-morpholin-4-
yl, CO-((4-loweralkyl)piperazin-1-yl), -NH(NH)NH², -NR⁴R^{4'} or lower alkyl,
with the proviso that a carbon atom is attached at the most to one heteroatom in
case this carbon atom is sp³-hybridized;

- 30 R⁴ and R^{4'} independently represent hydrogen; lower alkyl; cycloalkyl; cycloalkyl -
lower alkyl; hydroxy - lower alkyl; -COOR²; -CONH₂;

m and n represent the integer 0 or 1, with the proviso that in case m represents the integer 1, n is the integer 0, and in case n represents the integer 1, m is the integer 0;

- 5 p is the integer 1, 2, 3 or 4;
- r is the integer 3, 4, 5, or 6;
- s is the integer 2, 3, 4, or 5;
- t is the integer 1, 2, 3, or 4;
- u is the integer 1, 2, or 3;
- 10 v is the integer 2, 3, or 4;

and optically pure enantiomers, mixtures of enantiomers such as racemates, diastereomers, mixtures of diastereomers, diastereomeric racemates, mixtures of diastereomeric racemates, and the meso-form; as well as pharmaceutically acceptable salts, solvent complexes and morphological forms.

In the definitions of general formula I – if not otherwise stated – the term **lower alkyl**, alone or in combination with other groups, means saturated, straight and branched chain groups with one to seven carbon atoms, preferably one to four carbon atoms that can be optionally substituted by halogens. Examples of lower alkyl groups are methyl, ethyl, n-propyl, iso-propyl, n-butyl, iso-butyl, sec-butyl, tert-butyl, pentyl, hexyl and heptyl. The methyl, ethyl nad isopropyl groups are preferred.

25 The term **lower alkoxy** refers to a R-O group, wherein R is a lower alkyl. Examples of lower alkoxy groups are methoxy, ethoxy, propoxy, iso-propoxy, iso-butoxy, sec-butoxy and tert-butoxy.

The term **lower alkenyl**, alone or in combination with other groups, means
30 straight and branched chain groups comprising an olefinic bond and consisting of two to seven carbon atoms, preferably two to four carbon atoms, that can be

- optionally substituted by halogens. Examples of lower alkenyl are vinyl, propenyl or butenyl.
- 5 The term **lower alkynyl**, alone or in combination with other groups, means straight and branched chain groups comprising a triple bond and consisting of two to seven carbon atoms, preferably two to four carbon atoms, that can be optionally substituted by halogens. Examples of lower alkynyl are ethynyl, propynyl or butynyl.
- 10 The term **lower alkylene**, alone or in combination with other groups, means straight and branched divalent chain groups with one to seven carbon atoms, preferably one to four carbon atoms, that can be optionally substituted by halogens. Examples of lower alkylene are ethylene, propylene or butylene.
- 15 The term **lower alkenylene**, alone or in combination with other groups, means straight and branched divalent chain groups comprising an olefinic bond and consisting of two to seven carbon atoms, preferably two to four carbon atoms, that can be optionally substituted by halogens. Examples of lower alkenylene are vinylene, propenylene and butenylene.
- 20 The term **lower alkylenedioxy**, refers to a lower alkylene substituted at each end by an oxygen atom. Examples of lower alkylenedioxy groups are preferably methylenedioxy and ethylenedioxy.
- 25 The term **lower alkyleneoxy** refers to a lower alkylene substituted at one end by an oxygen atom. Examples of lower alkyleneoxy groups are preferably methyleneoxy, ethyleneoxy and propyleneoxy.
- 30 The term **halogen** means fluorine, chlorine, bromine or iodine, preferably fluorine, chlorine and bromine.

The term **cycloalkyl** alone or in combination, means a saturated cyclic hydrocarbon ring system with 3 to 7 carbon atoms, e.g. cyclopropyl, cyclobutyl, cyclopentyl, cyclohexyl and cycloheptyl, which can be optionally mono- or multisubstituted by lower alkyl, lower alkenyl, lower alkenylene, lower alkoxy, 5 lower alkylenoxy, lower alkylenedioxy, hydroxy, halogen, -CF₃, -NR¹R¹, -NR¹C(O)R¹, -NR¹S(O₂)R¹, -C(O)NR¹R¹, lower alkylcarbonyl, -COOR¹, -SR¹, -SOR¹, -SO₂R¹, -SO₂NR¹R¹ whereby R¹ represents hydrogen; lower alkyl; lower alkenyl; lower alkinyl; cycloalkyl; aryl; cycloalkyl - lower alkyl. The cyclopropyl group is a preferred group.

10

The term **aryl**, alone or in combination, relates to the phenyl, the naphthyl or the indanyl group, preferably the phenyl group, which can be optionally mono- or multisubstituted by lower alkyl, lower alkenyl, lower alkinyl, lower alkenylene or lower alkylene forming with the aryl ring a five- or six-membered ring, lower 15 alkoxy, lower alkylenedioxy, lower alkylenoxy, hydroxy, hydroxy-lower alkyl, halogen, cyano, -CF₃, -OCF₃, -NR¹R¹, -NR¹R¹ - lower alkyl, -NR¹C(O)R¹, -NR¹S(O₂)R¹, -C(O)NR¹R¹, -NO₂, lower alkylcarbonyl, -COOR¹, -SR¹, -SOR¹, -SO₂R¹, -SO₂NR¹R¹, benzyloxy, whereby R¹ has the meaning given above. Preferred substituents are halogen, lower alkoxy, lower alkyl, CF₃, OCF₃.

20

The term **aryloxy** refers to an Ar-O group, wherein Ar is an aryl. An example of a lower aryloxy group is phenoxy.

The term **heterocyclyl**, alone or in combination, means saturated or unsaturated 25 (but not aromatic) five-, six- or seven-membered rings containing one or two nitrogen, oxygen or sulfur atoms which may be the same or different and which rings can be optionally substituted with lower alkyl, hydroxy, lower alkoxy and halogen. The nitrogen atoms, if present, can be substituted by a -COOR² group. Examples of such rings are piperidinyl, morpholinyl, thiomorpholinyl, 30 piperazinyl, tetrahydropyranyl, dihydropyranyl, 1,4-dioxanyl, pyrrolidinyl, tetrahydrofuranyl, dihydropyrrolyl, imidazolidinyl, dihydropyrazolyl, pyrazolidinyl, dihydroquinolinyl, tetrahydroquinolinyl, tetrahydroisoquinolinyl.

The term **heteroaryl**, alone or in combination, means six-membered aromatic rings containing one to four nitrogen atoms; benzofused six-membered aromatic rings containing one to three nitrogen atoms; five-membered aromatic rings containing one oxygen, one nitrogen or one sulfur atom; benzofused five-membered aromatic rings containing one oxygen, one nitrogen or one sulfur atom; five-membered aromatic rings containing one oxygen and one nitrogen atom and benzofused derivatives thereof; five-membered aromatic rings containing a sulfur and a nitrogen or an oxygen atom and benzofused derivatives thereof; five-membered aromatic rings containing two nitrogen atoms and benzofused derivatives thereof; five-membered aromatic rings containing three nitrogen atoms and benzofused derivatives thereof, or a tetrazolyl ring. Examples of such ring systems are furanyl, thiophenyl, pyrrolyl, pyridinyl, pyrimidinyl, indolyl, quinolinyl, isoquinolinyl, imidazolyl, triazinyl, thiazinyl, thiazolyl, isothiazolyl, pyridazinyl, pyrazolyl, oxazolyl, isoxazolyl, coumarinyl, benzothiophenyl, quinazolinyl, quinoxalinyl. Such rings may be adequately substituted with lower alkyl, lower alkenyl, lower alkinyl, lower alkylene, lower alkenylene, lower alkylenedioxy, lower alkyleneoxy, hydroxy-lower alkyl, lower alkoxy, hydroxy, halogen, cyano, -CF₃, -OCF₃, -NR¹R¹, -NR¹R¹' - lower alkyl, -N(R¹)COR¹, -N(R¹)SO₂R¹, -CONR¹R¹', -NO₂, lower alkylcarbonyl, -COOR¹, -SR¹, -SOR¹, -SO₂R¹, -SO₂NR¹R¹', another aryl, another heteroaryl or another heterocyclyl and the like, whereby R¹ has the meaning given above.

The term **heteroaryloxy** refers to a Het-O group, wherein Het is a heteroaryl.

25

The term **sp³-hybridized** refers to a carbom atom and means that this carbon atom forms four bonds to four substituents placed in a tetragonal fashion around this carbon atom.

30

The expression **pharmaceutically acceptable salts** encompasses either salts with inorganic acids or organic acids like hydrochloric or hydrobromic acid, sulfuric acid, phosphoric acid, citric acid, formic acid, acetic acid, maleic acid, tartaric

acid, benzoic acid, methanesulfonic acid, p-toluenesulfonic acid, and the like that are non toxic to living organisms or in case the compound of formula I is acidic in nature with an inorganic base like an alkali or earth alkali base, e.g. sodium hydroxide, potassium hydroxide, calcium hydroxide and the like.

5

The compounds of the general formula I can contain two or more asymmetric carbon atoms and may be prepared in form of optically pure enantiomers, mixtures of enantiomers such as racemates, diastereomers, mixtures of diastereomers, diastereomeric racemates, mixtures of diastereomeric racemates, 10 and the meso-form and pharmaceutically acceptable salts therof.

The present invention encompasses all these forms. Mixtures may be separated in a manner known *per se*, i.e. by column chromatography, thin layer chromatography, HPLC or crystallization.

15 A group of preferred compounds are compounds of general formula I wherein X, W, V, U, T, Q, L, and M are as defined in general formula I above and wherein

n is 0 and

m is 1.

20

Another group of preferred compounds of general formula I are those wherein X, W, V, U, T, Q, M, m, and n are as defined in general formula I above and

L represents H; -COR^{3''}; -COOR^{3''}; -CONR^{2''}R^{3''};

25

whereby R^{2''} and R^{3''} represent independently lower alkyl, lower cycloalkyl - lower alkyl, which lower alkyl and lower cycloalkyl - lower alkyl groups are unsubstituted or monosubstituted with halogen, cyano, hydroxy, -OCOCH₃, -CONH₂, -COOH, -NH₂, with the proviso that a carbon atom is attached at the 30 most to one heteroatom in case this carbon atom is sp³-hybridized.

Another group of preferred compounds of general formula I above are those wherein X, W, V, U, L, m, and n are as defined in general formula I and

- T is -CONR¹;
- 5 Q is methylene;
- M is hydrogen; aryl; or heteroaryl.

Another group of even more preferred compounds of general formula I are those wherein X, W, U, L, T, Q, M, m, and n are as defined in general formula I above
10 and

V is -CH₂CH₂O-; -CH₂CH₂CH₂O-; -OCH₂CH₂O-; -.

Another group of also more preferred compounds of general formula I are those
15 wherein V, U, T, Q, M, L, m, and n are as defined in general formula I above and

X and W represent -CH-.

Another group of also more preferred compounds of general formula I are those
20 wherein X, W, V, Q, T, M, L, m, and n are as defined in general formula I above and

U is a mono-, di-, or trisubstituted phenyl wherein the substituents are halogen;
lower alkyl or lower alkoxy.

25 Especially preferred compounds of general formula I are those selected from the group consisting of:

(*rac.*)-(IR*, 5S*)-3-acetyl-7-{4-[2-(2,3-dichlorophenoxy)ethyl]phenyl}-3,9-di-
30 azabicyclo[3.3.1]non-6-ene-6-carboxylic acid methylphenethylamide;

(*rac.*)-(1*R*^{*}, 5*S*^{*})-3-acetyl-7-{4-[3-(2-chloro-5-methylphenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid methylphenethylamide;

5 (*rac.*)-(2-methoxyphenyl)acetic acid (1*R*^{*}, 5*S*^{*})-3-acetyl-7-{4-[3-(2-chlorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-en-6-ylmethyl ester;

(*rac.*)-(1*R*^{*}, 5*S*^{*})-3-acetyl-7-{4-[2-(4-bromo-3-methylphenoxy)ethyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid methylphenethylamide;

10 (*rac.*)-(2-methoxyphenyl)acetic acid (1*R*^{*}, 5*S*^{*})-3-acetyl-7-{4-[3-(2-chlorophenoxy)propoxy]phenyl}-3,9-diazabicyclo[3.3.1]non-6-en-6-ylmethyl ester;

(*rac.*)-(1*R*^{*}, 5*S*^{*})-3-acetyl-7-{4-[3-(2,4-dimethylphenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid methylphenethylamide;

15 (*rac.*)-(1*R*^{*}, 5*S*^{*})-3-acetyl-7-{4-[3-(4-bromo-3-methylphenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid methylphenethylamide;

20 (*rac.*)-(1*R*^{*}, 5*S*^{*})-3-acetyl-7-{4-[2-(2-acetylphenoxy)ethyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid methylphenethylamide;

(*rac.*)-(1*R*^{*}, 5*S*^{*})-3-acetyl-7-{4-[2-(2-ethylphenoxy)ethyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid methylphenethylamide;

25 (*rac.*)-(1*R*^{*}, 5*S*^{*})-3-acetyl-7-{4-[2-(3,5-dichlorophenoxy)ethyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid methylphenethylamide;

(*rac.*)-(1*R*^{*}, 5*S*^{*})-3-acetyl-7-{4-[3-(3,4-dichlorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid methylphenethylamide;

30 (*rac.*)-(1*R*^{*}, 5*S*^{*})-3-acetyl-7-{4-[2-(2,4-dimethylphenoxy)ethyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid methylphenethylamide;

- (*rac.*)-(1*R*^{*,} 5*S*^{*})-3-acetyl-7-[4-(3-*o*-tolyloxypropyl)phenyl]-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid methylphenethylamide;
- 5 (*rac.*)-(2-methoxyphenyl)acetic acid (1*R*^{*,} 5*S*^{*})-3-acetyl-7-{6-[3-(2-methoxybenzyloxy)propoxy]pyridin-3-yl}-3,9-diazabicyclo[3.3.1]non-6-en-6-ylmethyl ester;
- 10 (*rac.*)-(1*R*^{*,} 5*S*^{*})-3-acetyl-7-{4-[2-(2-chlorophenoxy)ethoxy]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid methylphenethylamide;
- 15 (*rac.*)-(1*R*^{*,} 5*S*^{*})-3-Acetyl-7-[4-(2-*m*-tolyloxyethyl)phenyl]-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid methylphenethylamide;
- 20 (*rac.*)-(1*R*^{*,} 5*S*^{*})-3-acetyl-7-{4-[2-(3-methoxyphenoxy)ethyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid methylphenethylamide;
- 25 (*rac.*)-(1*R*^{*,} 5*S*^{*})-3-acetyl-7-{4-[2-(2-chloro-3-trifluoromethylphenoxy)ethyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid methylphenethylamide;
- (*rac.*)-(2-methoxyphenyl)acetic acid (1*R*^{*,} 5*S*^{*})-3-acetyl-7-{4-[3-(2-bromophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-en-6-ylmethyl ester;
- 30 (*rac.*)-(1*R*^{*,} 5*S*^{*})-3-acetyl-7-{4-[3-(3-chlorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid methylphenethylamide;
- 35 (*rac.*)-(1*R*^{*,} 5*S*^{*})-3-acetyl-7-{4-[3-(2-isopropylphenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid methylphenethylamide;

- (*rac.*)-(1*R*^{*, 5*S*})-3-acetyl-7-{4-[2-(3-bromophenoxy)ethyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid methylphenethylamide;
- (*rac.*)-(1*R*S)-3-acetyl-7-{4-[3-(2-bromophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid methylphenethylamide;
- (*rac.*)-(2-methoxyphenyl)acetic acid (1*R*S)-3-acetyl-7-[4-(3-*o*-tolyloxypropyl)phenyl]-3,9-diazabicyclo[3.3.1]non-6-en-6-ylmethyl ester;
- 10 (*rac.*)-(1*R*S)-3-acetyl-7-{4-[2-(3,4-dichlorophenoxy)ethyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid methylphenethylamide;
- (*rac.*)-(1*R*S)-3-Acetyl-7-{4-[3-(2-*tert*-butylphenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid methylphenethylamide;
- 15 (*rac.*)-(1*R*S)-3-acetyl-7-{4-[2-(3-chlorophenoxy)ethyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid methylphenethylamide;
- (*rac.*)-(1*R*S)-3-acetyl-7-{4-[3-(2-ethylphenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid methylphenethylamide;
- 20 (*rac.*)-(1*R*S)-3-acetyl-7-{4-[3-(2,3-dimethylphenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid methylphenethylamide;
- (*rac.*)-(1*R*S)-3-acetyl-7-{4-[3-(2,3-dimethylphenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid methylphenethylamide;
- 25 (*rac.*)-(1*R*S)-3-acetyl-7-{4-[2-(3-trifluoromethoxyphenoxy)ethyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid methylphenethylamide;
- (*rac.*)-(1*R*S)-3-acetyl-7-{4-[3-(3,4-dimethylphenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid methylphenethylamide;
- 30 (*rac.*)-(1*R*S)-3-acetyl-7-{4-[3-(2-chloro-4,5-dimethylphenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid methylphenethylamide;

- (*rac.*)-(1*R*^{*,} 5*S*^{*})-3-acetyl-7-{4-[2-(2-chlorophenyl)ethoxy]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid methylphenethylamide;
- 5 (*rac.*)-(2-methoxyphenyl)acetic acid (1*R*^{*,} 5*S*^{*})-3-[2-(4-chlorophenyl)ethyl]-7-{4-[3-(2-methoxybenzyloxy)propoxy]phenyl}-3,9-diazabicyclo[3.3.1]non-6-en-6-yl methyl ester;
- 10 (*rac.*)-(1*R*^{*,} 5*S*^{*})-3-acetyl-7-{4-[3-(2-chloro-3-trifluoromethylphenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid methylphenethylamide;
- 15 (*rac.*)-(2-methoxyphenyl)acetic acid (1*R*^{*,} 5*S*^{*})-3-acetyl-7-{4-[2-(3-chlorophenoxy)ethyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-en-6-ylmethyl ester;
- (*rac.*)-(1*R*^{*,} 5*S*^{*})-3-acetyl-7-[4-(2-*m*-tolyloxyethoxy)phenyl]-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid methylphenethylamide;
- 20 (*rac.*)-(1*R*^{*,} 5*S*^{*})-3-acetyl-7-{4-[3-(2,4-dichlorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid methylphenethylamide;
- (*rac.*)-(1*R*^{*,} 5*S*^{*})-3-acetyl-7-{4-[3-(2,5-dimethylphenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid methylphenethylamide;
- 25 (*rac.*)-(1*R*^{*,} 5*S*^{*})-3-acetyl-7-{4-[2-(3,5-dimethylphenoxy)ethyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid methylphenethylamide;
- (*rac.*)-(1*R*^{*,} 5*S*^{*})-3-acetyl-7-{4-[2-(2-ethoxyphenoxy)ethyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid methylphenethylamide;
- 30 (*rac.*)-(2-methoxyphenyl)acetic acid (1*R*^{*,} 5*S*^{*})-3-acetyl-7-{4-[3-(3-chlorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-en-6-ylmethyl ester;

- (*rac.*)-(1*R*^{*,} 5*S*^{*})-3-acetyl-7-{4-[3-(3,5-dimethylphenoxy)propyl]phenyl}-3,9-di-azabicyclo[3.3.1]non-6-ene-6-carboxylic acid methylphenethylamide;
- 5 (*rac.*)-(1*R*^{*,} 5*S*^{*})-3-acetyl-7-{4-[3-(2-fluorophenoxy)propyl]phenyl}-3,9-diazabi-cyclo[3.3.1]non-6-ene-6-carboxylic acid methylphenethylamide;
- (*rac.*)-(2-methoxyphenyl)acetic acid (*1R*^{*,} 5*S*^{*})-3-acetyl-7-{4-[2-(2-chlorophen-oxy)ethyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-en-6-ylmethyl ester;
- 10 (*rac.*)-(1*R*^{*,} 5*S*^{*})-3-acetyl-7-{4-[2-(2,5-dichlorophenoxy)ethyl]phenyl}-3,9-diaza-bicyclo[3.3.1]non-6-ene-6-carboxylic acid methylphenethylamide;
- (*rac.*)-(1*R*^{*,} 5*S*^{*})-3-acetyl-7-{4-[3-(2,3,5-trimethylphenoxy)propyl]-phenyl}-3,9-di-azabicyclo[3.3.1]non-6-ene-6-carboxylic acid methylphenethylamide;
- 15 (*rac.*)-(1*R*^{*,} 5*S*^{*})-3-acetyl-7-{4-[2-(3-chlorophenyl)ethoxy]phenyl}-3,9-diazabi-cyclo[3.3.1]non-6-ene-6-carboxylic acid methylphenethylamide;
- (*rac.*)-(1*R*^{*,} 5*S*^{*})-3-acetyl-7-{4-[2-(3-chlorophenyl)ethoxy]phenyl}-3,9-diazabi-cyclo[3.3.1]non-6-ene-6-carboxylic acid methylphenethylamide;
- 20 (*rac.*)-(1*R*^{*,} 5*S*^{*})-3-acetyl-7-{4-[2-(2,3-difluorophenoxy)ethyl]phenyl}-3,9-diaza-bicyclo[3.3.1]non-6-ene-6-carboxylic acid methylphenethylamide;
- (*rac.*)-(1*R*^{*,} 5*S*^{*})-3-acetyl-7-{4-[2-(2-chlorophenoxy)ethyl]phenyl}-3,9-diazabi-cyclo[3.3.1]non-6-ene-6-carboxylic acid methylphenethylamide;
- 25 (*rac.*)-(1*R*^{*,} 5*S*^{*})-3-acetyl-7-{4-[2-(2,4-dichlorophenoxy)ethyl]phenyl}-3,9-diaza-bicyclo[3.3.1]non-6-ene-6-carboxylic acid methylphenethylamide;
- (*rac.*)-(1*R*^{*,} 5*S*^{*})-3-acetyl-7-{4-[2-(2,3,6-trifluorophenoxy)ethyl]phenyl}-3,9-di-azabicyclo[3.3.1]non-6-ene-6-carboxylic acid methylphenethylamide;

- (*rac.*)-(2-methoxyphenyl)acetic acid (*1R**, *5S**)-3-acetyl-7-{4-[3-(2-trifluoromethylphenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-en-6-ylmethyl ester;
- (*rac.*)-(*1R**, *5S**)-3-acetyl-7-{4-[2-(3-trifluoromethylphenoxy)ethyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid methylphenethylamide,
- (*rac.*)-(*1R**, *5S**)-3-acetyl-7-{4-[3-(3,5-dichlorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid methylphenethylamide;
- 10 (*rac.*)-(*1R**, *5S**)-3-acetyl-7-{4-[3-(*m*-tolyloxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid methylphenethylamide;
- (*rac.*)-(*1R**, *5S**)-3-acetyl-7-{4-[2-(2-methoxyphenyl)ethoxy]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid methylphenethylamide;
- 15 (*rac.*)-(*1R**, *5S**)-3-acetyl-7-{4-[3-(4-fluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid methylphenethylamide;
- (*rac.*)-(*1R**, *5S**)-3-acetyl-7-{4-[2-(3,4-dimethylphenoxy)ethyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid methylphenethylamide;
- 20 (*rac.*)-(2-methoxyphenyl)acetic acid (*1R**, *5S**)-3-acetyl-7-{4-[3-(3-fluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-en-6-ylmethyl ester;
- 25 (*rac.*)-(*1R**, *5S**)-3-acetyl-7-{4-[3-(2-methoxybenzyloxy)propoxy]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid phenethylamide;
- (*rac.*)-(2-methoxyphenyl)acetic acid (*1R**, *5S**)-3-acetyl-7-[4-(3-phenoxypropyl)phenyl]-3,9-diazabicyclo[3.3.1]non-6-en-6-ylmethyl ester;
- 30 (*rac.*)-(*1R**, *5S**)-3-acetyl-7-{4-[3-(2-acetyl-5-fluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid methylphenethylamide;

- (*rac.*)-(1*R*^{*}, 5*S*^{*})-7-{4-[3-(2-methoxybenzyloxy)propoxy]phenyl}-6-[2-(2-methoxyphenyl)acetoxymethyl]-3,9-diazabicyclo[3.3.1]non-6-ene-9-carboxylic acid 2,2,2-trichloro-1,1-dimethylethyl ester;
- 5 (*rac.*)-(1*R*^{*}, 5*S*^{*})-3-acetyl-7-{4-[3-(2-chlorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid methylphenethylamide;
- 10 (*rac.*)-(1*R*^{*}, 5*S*^{*})-3-acetyl-7-{4-[3-(2,3-difluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid methylphenethylamide;
- 15 (*rac.*)-(1*R*^{*}, 5*S*^{*})-3-acetyl-7-{4-[2-(2-*tert*-butylphenoxy)ethyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid methylphenethylamide;
- 20 1:1 mixture of (2*R*)- and (2*S*)-*N*-(1*R*^{*}, 5*S*^{*})-3-acetyl-7-{4-[3-(2-methoxybenzyl-oxy)propoxy]phenyl}-3,9-diazabicyclo[3.3.1]non-6-en-6-ylmethyl)-*N*-methyl-2-phenylpropionamide;
- 25 (*rac.*)-(1*R*^{*}, 5*S*^{*})-3-acetyl-7-{4-[3-(2-*tert*-butyl-4-methylphenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid methylphenethylamide;
- 30 (*rac.*)-(1*R*^{*}, 5*S*^{*})-3-acetyl-7-{4-[3-(2-*tert*-butyl-6-methylphenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid methylphenethylamide;
- (*rac.*)-(1*R*^{*}, 5*S*^{*})-3-acetyl-7-{4-[3-(4-chloro-2-methoxyphenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid methylphenethylamide;
- 35 (*rac.*)-(1*R*^{*}, 5*S*^{*})-3-acetyl-7-{4-[2-(2-methoxy-5-methylphenoxy)ethyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid methylphenethylamide;

- (*rac.*)-(1*R*^{*}, 5*S*^{*})-3-acetyl-7-{4-[3-(3-methoxyphenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid methylphenethylamide;
- 5 (*rac.*)-(1*R*^{*}, 5*S*^{*})-3-acetyl-7-{4-[3-(4-*tert*-butyl-2-methylphenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid methylphenethylamide;
- (rac.)-(1*R*^{*}, 5*S*^{*})-3-acetyl-7-{4-[2-(3,5-dimethoxyphenoxy)ethyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid methylphenethylamide;
- 10 (*rac.*)-(1*R*^{*}, 5*S*^{*})-3-acetyl-7-{4-[2-(2-*iso*-propylphenoxy)ethyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid methylphenethylamide;
- 15 (*rac.*)-(1*R*^{*}, 5*S*^{*})-3-acetyl-7-{4-[3-(2-methoxybenzyloxy)propoxy]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid [2-(2-methoxyphenyl)ethyl]amide;
- (rac.)-(1*R*^{*}, 5*S*^{*})-3-acetyl-7-{4-[3-(3-bromophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid methylphenethylamide;
- 20 (*rac.*)-(2-methoxyphenyl)acetic acid (*1R*^{*}, 5*S*^{*})-3-acetyl-7-[4-(3-*p*-tolyloxypropyl)phenyl]-3,9-diazabicyclo[3.3.1]non-6-en-6-ylmethyl ester;
- (rac.)-acetic acid 2-((1*R*^{*}, 5*S*^{*})-7-{4-[3-(2-chlorophenoxy)propyl]phenyl}-6-[(2-chlorophenyl)ethyl]methylcarbamoyl)-3,9-diazabicyclo[3.3.1]non-6-en-3-yl)-2-oxoethyl ester;
- 25 (*rac.*)-(1*R*^{*}, 5*S*^{*})-7-{4-[3-(2-chlorophenoxy)propyl]phenyl}-3-(2-cyanoacetyl)-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid [2-(2-chlorophenyl)ethyl]methylamide;
- 30

(*rac.*)-(1*R*^{*, 5*S*^{*})-3-(2-acetylaminoacetyl)-7-{4-[3-(2-chlorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid [2-(2-chlorophenyl)ethyl]methylamide;}

5 1:1 mixture of (*IR*, *5S*)-3-((4*R*)-2-acetylamino-4-methylpentanoyl)-7-{4-[3-(2-chlorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid [2-(2-chlorophenyl)ethyl]methylamide and (*IS*, *5R*)-3-((4*R*)-2-acetylamino-4-methylpentanoyl)-7-{4-[3-(2-chlorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid [2-(2-chlorophenyl)ethyl]methylamide;

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(*rac.*)-(1*R*^{*, 5*S*^{*})-3-acetyl-7-{4-[3-(2-chlorophenoxy)propyl]phenyl}-3,9-diazabi-cyclo[3.3.1]non-6-ene-6-carboxylic acid [2-(2-chlorophenyl)ethyl]methylamide;}

15

(*rac.*)-(1*R*^{*, 5*S*^{*})-3-acetyl-7-{4-[3-(2-chlorophenoxy)propyl]phenyl}-3,9-diazabi-cyclo[3.3.1]non-6-ene-6-carboxylic acid (2-hydroxybenzyl)methylamide;}

20

1:1 mixture of (*rac.*)-(1*R*^{*, 5*S*^{*})-3-acetyl-7-{4-[3-(2-chlorophenoxy)propyl]-phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid [(3*R*^{*})-3-(2-chloro-phenyl)butyl]methylamide and (*rac.*)-(1*R*^{*, 5*S*^{*})-3-acetyl-7-{4-[3-(2-chlorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid [(3*S*^{*})-3-(2-chlorophenyl)butyl]methylamide;}}

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(*rac.*)-(1*R*^{*, 5*S*^{*})-3-acetyl-7-{4-[3-(2-chlorophenoxy)propyl]phenyl}-3,9-diazabi-cyclo[3.3.1]non-6-ene-6-carboxylic acid methyl-(4-phenylbutyl)amide;}

(*rac.*)-(1*R*^{*, 5*S*^{*})-3-acetyl-7-{4-[3-(2-chlorophenoxy)propyl]phenyl}-3,9-diazabi-cyclo[3.3.1]non-6-ene-6-carboxylic acid methyl-(3-phenoxypropyl)amide;}

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(*rac.*)-(1*R*^{*, 5*S*^{*})-3-acetyl-7-{4-[3-(2-chlorophenoxy)propyl]phenyl}-3,9-diazabi-cyclo[3.3.1]non-6-ene-6-carboxylic acid methyl-(4-phenylpentyl)amide;}

- (*rac.*)-(1*R*^{*,} 5*S*^{*})-3-acetyl-7-{4-[3-(2-chlorophenoxy)propyl]phenyl}-3,9-diazabi-cyclo[3.3.1]non-6-ene-6-carboxylic acid (3-benzo[1,3]dioxol-5-ylpropyl)methylamide;

5 (*rac.*)-(1*R*^{*,} 5*S*^{*})-3-acetyl-7-{4-[3-(2-chlorophenoxy)propyl]phenyl}-3,9-diazabi-cyclo[3.3.1]non-6-ene-6-carboxylic acid [2-(4-methoxyphenoxy)ethyl]methylamide;

10 (*rac.*)-(1*R*^{*,} 5*S*^{*})-3-acetyl-7-{4-[3-(2-chlorophenoxy)propyl]phenyl}-3,9-diazabi-cyclo[3.3.1]non-6-ene-6-carboxylic acid [2-(4-chlorophenoxy)ethyl]methylamide;

15 (*rac.*)-(1*R*^{*,} 5*S*^{*})-3-acetyl-7-{4-[3-(2-chlorophenoxy)propyl]phenyl}-3,9-diazabi-cyclo[3.3.1]non-6-ene-6-carboxylic acid methyl-(2-*p*-tolyloxyethyl)amide;

20 (*rac.*)-(1*R*^{*,} 5*S*^{*})-3-acetyl-7-{4-[3-(2-chlorophenoxy)propyl]phenyl}-3,9-diazabi-cyclo[3.3.1]non-6-ene-6-carboxylic acid diethylamide;

25 (*rac.*)-(1*R*^{*,} 5*S*^{*})-3-acetyl-7-{4-[3-(2-bromo-5-fluorophenoxy)ethyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid (2-chlorobenzyl)cyclopropylamide;

30 (*rac.*)-(1*R*^{*,} 5*S*^{*})-3-acetyl-7-{4-[3-(2-chlorophenoxy)propyl]phenyl}-3,9-diaza-bicyclo[3.3.1]non-6-ene-6-carboxylic acid (2-chlorobenzyl)cyclopropylamide;

35 (*rac.*)-(1*R*^{*,} 5*S*^{*})-3-acetyl-7-{4-[3-(2-chlorophenoxy)propyl]phenyl}-3,9-diaza-bicyclo[3.3.1]non-6-ene-6-carboxylic acid [2-(4-methoxyphenoxy)ethyl]methyl-

- (*rac.*)-(1*R*^{*}, 5*S*^{*})-3-acetyl-7-{4-[3-(2-chlorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid methyl-(3-trifluoromethylbenzyl)amide;
- 5 (*rac.*)-(1*R*^{*}, 5*S*^{*})-3-acetyl-7-{4-[3-(2-chlorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid [2-(3,4-dimethylphenoxy)ethyl]methylamide;
- 10 (*rac.*)-(1*R*^{*}, 5*S*^{*})-3-acetyl-7-{4-[3-(2-chlorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid (3,5-dimethoxybenzyl)methylamide;
- (*rac.*)-(1*R*^{*}, 5*S*^{*})-3-acetyl-7-{4-[3-(2-chlorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid (2-chlorobenzyl)methylamide;
- 15 (*rac.*)-(1*R*^{*}, 5*S*^{*})-3-acetyl-7-{4-[2-(2-chloro-4,5-dimethylphenoxy)ethoxy]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid methylphenethylamide;
- 20 (*rac.*)-(1*R*^{*}, 5*S*^{*})-3-acetyl-7-{4-[2-(3,4-dichlorophenoxy)ethoxy]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid methylphenethylamide;
- (*rac.*)-(1*R*^{*}, 5*S*^{*})-3-acetyl-7-{4-[3-(2-chlorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid (2-chlorobenzyl)ethylamide;
- 25 (*rac.*)-(1*R*^{*}, 5*S*^{*})-3-acetyl-7-{4-[3-(2,3,6-trichlorophenoxy)ethyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid methylphenethylamide;
- (*rac.*)-(1*R*^{*}, 5*S*^{*})-3-acetyl-7-{4-[3-(3-fluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid (2-chlorobenzyl)cyclopropyl-amide;
- 30 (*rac.*)-(1*R*^{*}, 5*S*^{*})-3-acetyl-7-{4-[3-(2-chlorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid (2-chlorobenzyl)isopropylamide;

(*rac.*)-(1*R*^{*, 5*S*})-5-[7-{4-[2-(2-bromo-5-fluorophenoxy)ethyl]phenyl}-6-(methylphenethylcarbamoyl)-3,9-diazabicyclo[3.3.1]non-6-en-3-yl]-2,2-dimethyl-5-oxo-pentanoic acid;

5

(1*S*, 5*R*)-3-acetyl-7-{4-[2-(2,3,6-trifluorophenoxy)ethyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid methylphenethylamide;

1:1-mixture of (*rac.*)-(1*R*S)-3-acetyl-7-{4-[3-(2-bromo-5-fluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid cyclopropyl-[2-((2*R*^{*})-2,3-dihydroxypropyl)benzyl]amide and (*rac.*)-(1*R*S)-3-acetyl-7-{4-[3-(2-bromo-5-fluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid cyclopropyl-[2-((2*S*^{*})-2,3-dihydroxypropyl)benzyl]amide;

15 (*rac.*)-(1*R*S)-3-acetyl-7-{4-[3-(2-bromo-5-fluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid cyclopropyl-[2-(2-hydroxyethyl)benzyl]amide;

20 (*rac.*)-(1*R*S)-3-acetyl-7-{4-[3-(2-bromo-5-fluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid benzylcyclopropylamide;

(*rac.*)-(1*R*S)-3-acetyl-7-{4-[3-(2-bromo-5-fluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid (2-chlorobenzyl)ethylamide;

25 (*rac.*)-(1*R*S)-3-acetyl-7-{4-[3-(2-bromo-5-fluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid cyclopropyl-(2-fluorobenzyl)amide;

30 (*rac.*)-(1*R*S)-3-acetyl-7-{4-[3-(2-bromo-5-fluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid cyclopropyl-(2-methylbenzyl)amide;

- (*rac.*)-(1*R*^{*}, 5*S*^{*})-3-acetyl-7-{4-[3-(2-bromo-5-fluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid cyclopropyl-[2-(4-methoxyphenoxy)ethyl]amide;
- 5 (*rac.*)-(1*R*^{*}, 5*S*^{*})-3-acetyl-7-{4-[3-(2-bromo-5-fluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid cyclopropyl-(2-*m*-tolyloxyethyl)amide;
- 10 (*rac.*)-(1*R*^{*}, 5*S*^{*})-3-acetyl-7-{4-[3-(2-bromo-5-fluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid cyclopropyl-[2-(3,4-dimethylphenoxy)ethyl]amide;
- 15 (*rac.*)-(1*R*^{*}, 5*S*^{*})-3-acetyl-7-{4-[3-(2-bromo-5-fluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid cyclopropylphenethylamide;
- 20 (*rac.*)-(1*R*^{*}, 5*S*^{*})-3-acetyl-7-{4-[3-(2-bromo-5-fluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid [2-(2-chlorophenyl)ethyl]-cyclopropylamide;
- 25 (*rac.*)-(1*R*^{*}, 5*S*^{*})-3-acetyl-7-{4-[3-(2-bromo-5-fluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid cyclopropyl-[2-(2,3-difluorophenyl)ethyl]amide;
- 30 (*rac.*)-(1*R*^{*}, 5*S*^{*})-3-acetyl-7-{4-[3-(2-bromo-5-fluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid cyclopropyl-[2-(4-fluorophenyl)ethyl]amide;
- 35 (*rac.*)-(1*R*^{*}, 5*S*^{*})-3-acetyl-7-{4-[3-(2-bromo-5-fluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid cyclopropyl-(2-*o*-tolylethyl)amide;

1:1-mixture of (*rac.*)-(1*R*^{*}, 5*S*^{*})-3-acetyl-7-{4-[3-(2-bromo-5-fluorophenoxy)-propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid ((2*R*^{*})-2-hydroxy-2-phenylethyl)methylamide and (*rac.*)-(1*R*^{*}, 5*S*^{*})-3-acetyl-7-{4-[3-(2-bromo-5-fluorophenoxy)-propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid ((2*S*^{*})-2-hydroxy-2-phenylethyl)methylamide;

(*rac.*)-(1*R*^{*}, 5*S*^{*})-3-acetyl-7-{4-[3-(2-bromo-5-fluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid cyclopropyl-(3-trifluoromethylbenzyl)amide;

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(*rac.*)-(1*R*^{*}, 5*S*^{*})-3-acetyl-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid cyclopropyl-(2-*o*-tolylethyl)amide;

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(*rac.*)-(1*R*^{*}, 5*S*^{*})-6-[(2-chlorobenzyl)ethylcarbamoyl]-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-9-carboxylic acid 2,2,2-trichloro-1,1-dimethylethyl ester;

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(*rac.*)-(1*R*^{*}, 5*S*^{*})-6-[(2-fluorobenzyl)cyclopropylcarbamoyl]-7-{4-[3-(2,3,6-trifluoro-phenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-9-carboxylic acid 2,2,2-trichloro-1,1-dimethylethyl ester;

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(*rac.*)-(1*R*^{*}, 5*S*^{*})-6-[(2-methylbenzyl)cyclopropylcarbamoyl]-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-9-carboxylic acid 2,2,2-trichloro-1,1-dimethylethyl ester;

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(*rac.*)-(1*R*^{*}, 5*S*^{*})-6-[cyclopropyl-(2-*o*-tolylethyl)carbamoyl]-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-9-carboxylic acid 2,2,2-trichloro-1,1-dimethylethyl ester;

- (*rac.*)-(1*R*^{*}, 5*S*^{*})-6-[cyclopropyl-(2-*p*-tolylethyl)carbamoyl]-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-9-carboxylic acid 2,2,2-trichloro-1,1-dimethylethyl ester;
- 5 (*rac.*)-(1*R*^{*}, 5*S*^{*})-6-[cyclopropyl-(3,5-dimethoxybenzyl)carbamoyl]-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-9-carboxylic acid 2,2,2-trichloro-1,1-dimethylethyl ester;
- 10 (*rac.*)-(1*R*^{*}, 5*S*^{*})-3-(4-carbamoylbutyryl)-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid cyclopropyl-[2-(4-methoxyphenoxy)ethyl]amide;
- 15 (*rac.*)-(1*R*^{*}, 5*S*^{*})-3-(4-carbamoylbutyryl)-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid cyclopropyl-[2-(3-methoxyphenoxy)ethyl]amide;
- 20 (*rac.*)-(1*R*^{*}, 5*S*^{*})-3-(4-carbamoylbutyryl)-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid cyclopropyl-(2-*m*-tolyloxyethyl)amide;
- 25 (*rac.*)-(1*R*^{*}, 5*S*^{*})-3-(4-carbamoylbutyryl)-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid cyclopropyl-[2-(3,4-dimethylphenoxy)ethyl]amide;
- 30 (*rac.*)-(1*R*^{*}, 5*S*^{*})-3-(4-carbamoylbutyryl)-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid cyclopropyl-[2-(4-fluorophenyl)ethyl]amide;

- (*rac.*)-(1*R*^{*}, 5*S*^{*})-3-(4-carbamoylbutyryl)-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid cyclopropyl-(2-*o*-tolylethyl)amide;
- 5 (*rac.*)-(1*R*^{*}, 5*S*^{*})-3-(4-carbamoylbutyryl)-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid cyclopropyl-(2-*p*-tolylethyl)amide;
- 10 1:1-mixture of (1*R*, 5*S*)-3-((1*S*, 4*R*)-4-hydroxypyrrolidine-2-carbonyl)-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid benzylcyclopropylamide and (1*S*, 5*R*)-3-((1*S*, 4*R*)-4-hydroxypyrrolidine-2-carbonyl)-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid benzylcyclopropylamide;
- 15 1:1-mixture of (1*R*, 5*S*)-3-((1*S*, 4*R*)-4-hydroxypyrrolidine-2-carbonyl)-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid (2-chlorobenzyl)ethylamide and (1*S*, 5*R*)-3-((1*S*, 4*R*)-4-hydroxypyrrolidine-2-carbonyl)-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid (2-chlorobenzyl)ethylamide;
- 20 1:1-mixture of (1*R*, 5*S*)-3-((1*S*, 4*R*)-4-hydroxypyrrolidine-2-carbonyl)-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid cyclopropyl-(2-fluorobenzyl)amide and (1*S*, 5*R*)-3-((1*S*, 4*R*)-4-hydroxypyrrolidine-2-carbonyl)-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo-[3.3.1]non-6-ene-6-carboxylic acid cyclopropyl-(2-fluorobenzyl)amide;
- 25 1:1-mixture of (1*R*, 5*S*)-3-((1*S*, 4*R*)-4-hydroxypyrrolidine-2-carbonyl)-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid cyclopropyl-(3-trifluoromethylbenzyl)amide and (1*S*, 5*R*)-3-((1*S*, 4*R*)-4-hydroxypyrrolidine-2-carbonyl)-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]-

- phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid cyclopropyl-(3-trifluoromethylbenzyl)amide;
- 1:1-mixture of (*IR, 5S*)-3-((*IS, 4R*)-4-hydroxypyrrolidine-2-carbonyl)-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid cyclopropyl-(2-methylbenzyl)amide and (*IS, 5R*)-3-((*IS, 4R*)-4-hydroxypyrrolidine-2-carbonyl)-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid cyclopropyl-(2-methylbenzyl)amide;
- 10 1:1-mixture of (*IR, 5S*)-3-((*IS, 4R*)-4-hydroxypyrrolidine-2-carbonyl)-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diaza-bicyclo[3.3.1]non-6-ene-6-carboxylic acid cyclopropyl-(2-*m*-tolyloxyethyl)amide and (*IS, 5R*)-3-((*IR 4S*)-4-hydroxypyrrolidine-2-carbonyl)-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid cyclopropyl-(2-*m*-tolyloxyethyl)amide;
- 15 1:1-mixture of (*IR, 5S*)-3-((*IS, 4R*)-4-hydroxypyrrolidine-2-carbonyl)-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid cyclopropyl-(3,5-dimethoxybenzyl)amide and (*IS, 5R*)-3-((*IS, 4R*)-4-hydroxypyrrolidine-2-carbonyl)-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid cyclopropyl-(3,5-dimethoxybenzyl)amide;
- 20 25 1:1-mixture of (*IR, 5S*)-3-((*IS, 4R*)-hydroxypyrrolidine-2-carbonyl)-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid cyclopropyl-(2-*p*-tolylethyl)amide and (*IS, 5R*)-3-((*IS, 4R*)-hydroxypyrrolidine-2-carbonyl)-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid cyclopropyl-(2-*p*-tolylethyl)-
- 30 amide;

- 1:1-mixture of (*rac.*)-(3*R**)-5-((1*R**, 5*S**)-6-{cyclopropyl-[2-(3-methoxyphenoxy)ethyl]carbamoyl}-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-en-3-yl)-3-hydroxy-5-oxopentanoic acid and (*rac.*)-(3*S**)-5-((1*R**, 5*S**)-6-{cyclopropyl-[2-(3-methoxyphenoxy)ethyl]carbamoyl}-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-en-3-yl)-3-hydroxy-5-oxopentanoic acid;
- 1:1-mixture of (*rac.*)-(3*R**)-5-((1*R**, 5*S**)-6-{cyclopropyl-[2-(3,4-dimethylphenoxy)ethyl]carbamoyl}-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-en-3-yl)-3-hydroxy-5-oxopentanoic acid and (*rac.*)-(3*S**)-5-((1*R**, 5*S**)-6-{cyclopropyl-[2-(3,4-dimethylphenoxy)ethyl]carbamoyl}-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-en-3-yl)-3-hydroxy-5-oxopentanoic acid;
- 15 1:1-mixture of (*rac.*)-(3*R**)-5-((1*R**, 5*S**)-(6-(cyclopropylphenethyl-carbamoyl)-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-en-3-yl)-3-hydroxy-5-oxopentanoic acid and (*rac.*)-(3*S**)-5-((1*R**, 5*S**)-5-(6-(cyclopropylphenethylcarbamoyl)-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-en-3-yl)-3-hydroxy-5-oxopentanoic acid;
- 20 1:1-mixture of (*rac.*)-(3*R**)-5-((1*R**, 5*S**)-6-{cyclopropyl-[2-(2,3-difluorophenyl)ethyl]carbamoyl}-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-en-3-yl)-3-hydroxy-5-oxopentanoic acid and (*rac.*)-(3*S**)-5-((1*R**, 5*S**)-6-{cyclopropyl-[2-(2,3-difluorophenyl)ethyl]carbamoyl}-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-en-3-yl)-3-hydroxy-5-oxopentanoic acid;
- 25 1:1-mixture of (*rac.*)-(3*R**)-5-((1*R**, 5*S**)-6-{cyclopropyl-[2-(4-fluorophenyl)ethyl]carbamoyl}-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-en-3-yl)-3-hydroxy-5-oxopentanoic acid and (*rac.*)-(3*S**)-5-((1*R**, 5*S**)-6-{cyclopropyl-[2-(4-fluorophenyl)ethyl]carbamoyl}-7-{4-[3-(2,3,6-

trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-en-3-yl)-3-hydroxy-5-oxopentanoic acid;

1:1-mixture of (*rac.*)-(3*R**)-5-((*IR**, 5*S**)-6-[cyclopropyl-(2-*o*-tolylethyl)-carbamoyl]-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-en-3-yl)-3-hydroxy-5-oxopentanoic acid and (*rac.*)-(3*S**)-5-((*IR**, 5*S**)-6-[cyclopropyl-(2-*o*-tolylethyl)carbamoyl]-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-en-3-yl)-3-hydroxy-5-oxopentanoic acid;

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1:1-mixture of (*rac.*)-(3*R**)-5-((*IR**, 5*S**)-6-[cyclopropyl-(2-*p*-tolylethyl)-carbamoyl]-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-en-3-yl)-3-hydroxy-5-oxopentanoic acid and (*rac.*)-(3*R**)-5-((*IR**, 5*S**)-6-[cyclopropyl-(2-*p*-tolylethyl)carbamoyl]-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-en-3-yl)-3-hydroxy-5-oxopentanoic acid;

(*rac.*)-(*IR**, 5*S**)-3-acetyl-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid benzylcyclopropylamide;

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(*rac.*)-(*IR**, 5*S**)-3-acetyl-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid cyclopropyl-(2-fluorobenzyl)-amide;

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(*rac.*)-(*IR**, 5*S**)-3-acetyl-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid cyclopropyl-(2-methylbenzyl)-amide;

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(*rac.*)-(*IR**, 5*S**)-3-acetyl-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid cyclopropyl-[2-(4-methoxyphenoxy)ethyl]amide;

- (*rac.*)-(1*R*^{*,} 5*S*^{*})-3-acetyl-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid cyclopropyl-[2-(3-methoxyphenoxy)ethyl]amide;
- 5 (*rac.*)-(1*R*^{*,} 5*S*^{*})-3-acetyl-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid cyclopropyl-[2-(3-methylphenoxy)ethyl]amide;
- (*rac.*)-(1*R*^{*,} 5*S*^{*})-3-acetyl-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid cyclopropylphenethylamide;
- 10 (*rac.*)-(1*R*^{*,} 5*S*^{*})-3-acetyl-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid cyclopropyl(2-*p*-tolylethyl)amide;
- 15 (*rac.*)-5-((1*R*^{*,} 5*S*^{*})-6-{cyclopropyl-[2-(4-methoxyphenoxy)ethyl]carbamoyl}-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-en-3-yl)-5-oxopentanoic acid;
- (*rac.*)-5-((1*R*^{*,} 5*S*^{*})-6-{cyclopropyl-[2-(3-methoxyphenoxy)ethyl]carbamoyl}-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-en-3-yl)-5-oxopentanoic acid;
- 20 (*rac.*)-5-((1*R*^{*,} 5*S*^{*})-6-{cyclopropyl[2-(3-methylphenoxy)ethyl]carbamoyl}-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-en-3-yl)-5-oxopentanoic acid;
- 25 (*rac.*)-5-((1*R*^{*,} 5*S*^{*})-6-{cyclopropyl[2-(3-methylphenoxy)ethyl]carbamoyl}-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-en-3-yl)-5-oxopentanoic acid;
- (*rac.*)-5-((1*R*^{*,} 5*S*^{*})-6-{cyclopropyl-[2-(3,4-dimethylphenoxy)ethyl]carbamoyl}-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-en-3-yl)-5-oxopentanoic acid;

- (*rac.*)-5-((*1R*^{*}, *5S*^{*})-6-(cyclopropylphenethylcarbamoyl)-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-en-3-yl)-5-oxopentanoic acid;
- 5 (*rac.*)-5-((*1R*^{*}, *5S*^{*})-6-{cyclopropyl-[2-(2,3-difluorophenyl)ethyl]carbamoyl}-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-en-3-yl)-5-oxopentanoic acid;
- 10 (*rac.*)-5-((*1R*^{*}, *5S*^{*})-6-{cyclopropyl-[2-(4-fluorophenyl)ethyl]carbamoyl}-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-en-3-yl)-5-oxopentanoic acid;
- 15 (*rac.*)-5-((*1R*^{*}, *5S*^{*})-6-{cyclopropyl-[2-(2-methylphenyl)ethyl]carbamoyl}-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-en-3-yl)-5-oxopentanoic acid;
- 20 (*rac.*)-5-((*1R*^{*}, *5S*^{*})-6-(benzylcyclopropylcarbamoyl)-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-en-3-yl)-5-oxopentanoic acid methyl ester;
- 25 (*rac.*)-5-((*1R*^{*}, *5S*^{*})-6-[cyclopropyl-(3-trifluoromethylbenzyl)carbamoyl]-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-en-3-yl)-5-oxopentanoic acid methyl ester;
- 30 (*rac.*)-5-((*1R*^{*}, *5S*^{*})-6-[cyclopropyl-(2-methylmethylbenzyl)carbamoyl]-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-en-3-yl)-5-oxopentanoic acid methyl ester;
- 35 (*rac.*)-5-((*1R*^{*}, *5S*^{*})-6-{cyclopropyl-[2-(4-methoxyphenoxy)ethyl]carbamoyl}-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-en-3-yl)-5-oxopentanoic acid methyl ester;

- (*rac.*.)-5-((*1R**, *5S**)-6-{cyclopropyl-[2-(3-methoxyphenoxy)ethyl]carbamoyl}-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-en-3-yl)-5-oxopentanoic acid methyl ester;
- 5 (*rac.*.)-5-((*1R**, *5S**)-6-{cyclopropyl-[2-(3-methylphenoxy)ethyl]carbamoyl}-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-en-3-yl)-5-oxopentanoic acid methyl ester;
- 10 (*rac.*.)-5-((*1R**, *5S**)-6-{cyclopropyl-[2-(3,4-dimethylphenoxy)ethyl]carbamoyl}-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-en-3-yl)-5-oxopentanoic acid methyl ester;
- 15 (*rac.*.)-5-((*1R**, *5S**)-6-(cyclopropylphenethylcarbamoyl)-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-en-3-yl)-5-oxopentanoic acid methyl ester;
- 20 (*rac.*.)-5-((*1R**, *5S**)-6-{{[2-(2-chlorophenyl)ethyl]cyclopropylcarbamoyl}-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-en-3-yl)-5-oxopentanoic acid methyl ester;
- 25 (*rac.*.)-5-((*1R**, *5S**)-6-{cyclopropyl-[2-(2,3-difluorophenyl)ethyl]carbamoyl}-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-en-3-yl)-5-oxopentanoic acid methyl ester;
- 30 (*rac.*.)-5-((*1R**, *5S**)-6-{cyclopropyl-[2-(4-fluorophenyl)ethyl]carbamoyl}-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-en-3-yl)-5-oxopentanoic acid methyl ester;
- 35 (*rac.*.)-5-((*1R**, *5S**)-6-{cyclopropyl-[2-(2-methylphenyl)ethyl]carbamoyl}-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-en-3-yl)-5-oxopentanoic acid methyl ester;

- (*rac.*)-5-((*1R*^{*}, *5S*^{*})-6-{cyclopropyl-[2-(4-methylphenyl)ethyl]carbamoyl}-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-en-3-yl)-5-oxopentanoic acid methyl ester;
- 5 (*rac.*)-5-((*1R*^{*}, *5S*^{*})-6-(benzylcyclopropylcarbamoyl)-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-en-3-yl)-2,2-dimethyl-5-oxopentanoic acid;
- 10 (*rac.*)-5-((*1R*^{*}, *5S*^{*})-6-[cyclopropyl-(3-trifluoromethylbenzyl)carbamoyl]-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-en-3-yl)-2,2-dimethyl-5-oxopentanoic acid;
- 15 (*rac.*)-5-((*1R*^{*}, *5S*^{*})-6-{cyclopropyl-[2-(4-methoxyphenoxy)ethyl]carbamoyl}-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-en-3-yl)-2,2-dimethyl-5-oxopentanoic acid;
- 20 (*rac.*)-5-((*1R*^{*}, *5S*^{*})-6-{cyclopropyl-[2-(3-methoxyphenoxy)ethyl]carbamoyl}-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-en-3-yl)-2,2-dimethyl-5-oxopentanoic acid;
- 25 (*rac.*)-5-((*1R*^{*}, *5S*^{*})-6-{cyclopropyl-[2-(3-methylphenoxy)ethyl]carbamoyl}-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-en-3-yl)-2,2-dimethyl-5-oxopentanoic acid;
- 30 (*rac.*)-5-((*1R*^{*}, *5S*^{*})-6-(cyclopropylphenethylcarbamoyl)-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-en-3-yl)-2,2-dimethyl-5-oxo-pentanoic acid;

- (*rac.*)-5-((*1R*^{*,} *5S*^{*})-6-{[2-(2-chlorophenyl)ethyl]cyclopropylcarbamoyl}-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-en-3-yl)-2,2-dimethyl-5-oxopentanoic acid;
- 5 (*rac.*)-5-((*1R*^{*,} *5S*^{*})-6-{[2-(2,3-difluorophenyl)ethyl]cyclopropylcarbamoyl}-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-en-3-yl)-2,2-dimethyl-5-oxopentanoic acid;
- 10 (*rac.*)-5-((*1R*^{*,} *5S*^{*})-6-{[2-(4-fluorophenyl)ethyl]cyclopropylcarbamoyl}-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-en-3-yl)-2,2-dimethyl-5-oxopentanoic acid;
- 15 (*rac.*)-5-((*1R*^{*,} *5S*^{*})-6-{[2-(2-methylphenyl)ethyl]cyclopropylcarbamoyl}-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-en-3-yl)-2,2-dimethyl-5-oxopentanoic acid;
- 20 (*rac.*)-5-((*1R*^{*,} *5S*^{*})-6-{[2-(4-methylphenyl)ethyl]cyclopropylcarbamoyl}-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-en-3-yl)-2,2-dimethyl-5-oxopentanoic acid;
- 25 1:1-mixture of (*rac.*)-(2*R*^{*,} 3*S*^{*})-4-((*1R*^{*,} *5S*^{*})-6-[(2-chlorobenzyl)ethylcarbamoyl]-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-en-3-yl)-2,3-dihydroxy-4-oxobutyric acid and (*rac.*)-(2*S*^{*,} 3*R*^{*})-4-((*1R*^{*,} *5S*^{*})-6-[(2-chlorobenzyl)ethylcarbamoyl]-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-en-3-yl)-2,3-dihydroxy-4-oxobutyric acid;
- 30 1:1-mixture of (*rac.*)-(2*R*^{*,} 3*S*^{*})-4-((*1R*^{*,} *5S*^{*})-6-[cyclopropyl-(2-fluorobenzyl)carbamoyl]-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-en-3-yl)-2,3-dihydroxy-4-oxobutyric acid and (*rac.*)-(2*S*^{*,} 3*R*^{*})-4-((*1R*^{*,} *5S*^{*})-6-[cyclopropyl-(2-fluorobenzyl)carbamoyl]-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-en-3-yl)-2,3-dihydroxy-4-oxobutyric acid;

phenoxy)propyl]phenyl}-3,9-diazabicyclo-[3.3.1]non-6-en-3-yl)-2,3-dihydroxy-4-oxobutyric acid;

- 5 1:1-mixture of (*rac.*)-(2*R*^{*,} 3*S*^{*})-4-((1*R*^{*,} 5*S*^{*})-6-[cyclopropyl-(3-trifluoromethylbenzyl)carbamoyl]-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-en-3-yl)-2,3-dihydroxy-4-oxobutyric acid and (*rac.*)-(2*S*[,] 3*R*^{*})-4-((1*R*^{*,} 5*S*^{*})-6-[cyclopropyl-(3-trifluoromethylbenzyl)carbamoyl]-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo-[3.3.1]non-6-en-3-yl)-2,3-dihydroxy-4-oxobutyric acid;
- 10 1:1-mixture of (*rac.*)-(2*R*[,] 3*S*^{*})-4-((1*R*[,] 5*S*^{*})-6-{cyclopropyl-[2-(2,3-difluorophenyl)ethyl]carbamoyl}-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-en-3-yl)-2,3-dihydroxy-4-oxobutyric acid and (*rac.*)-(2*S*[,] 3*R*^{*})-4-((1*R*[,] 5*S*^{*})-6-{cyclopropyl-[2-(2,3-difluorophenyl)ethyl]carbamoyl}-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo-[3.3.1]non-6-en-3-yl)-2,3-dihydroxy-4-oxobutyric acid;
- 15 1:1-mixture of (*rac.*)-(2*R*[,] 3*S*^{*})-4-((1*R*[,] 5*S*^{*})-6-{cyclopropyl-[2-(2-methylphenyl)ethyl]carbamoyl}-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-en-3-yl)-2,3-dihydroxy-4-oxobutyric acid and (*rac.*)-(2*S*[,] 3*R*^{*})-4-((1*R*[,] 5*S*^{*})-6-{cyclopropyl-[2-(2-methylphenyl)ethyl]carbamoyl}-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo-[3.3.1]non-6-en-3-yl)-2,3-dihydroxy-4-oxobutyric acid;
- 20 1:1-mixture of (*rac.*)-(2*R*[,] 3*S*^{*})-4-((1*R*[,] 5*S*^{*})-6-[cyclopropyl-(3,5-dimethoxybenzyl)carbamoyl]-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-en-3-yl)-2,3-dihydroxy-4-oxobutyric acid and (*rac.*)-(2*S*[,] 3*R*^{*})-4-((1*R*[,] 5*S*^{*})-6-[cyclopropyl-(3,5-dimethoxybenzyl)carbamoyl]-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-en-3-yl)-2,3-dihydroxy-4-oxobutyric acid;
- 25 1:1-mixture of (*rac.*)-(2*R*[,] 3*S*^{*})-4-((1*R*[,] 5*S*^{*})-6-[cyclopropyl-(3,5-dimethoxybenzyl)carbamoyl]-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-en-3-yl)-2,3-dihydroxy-4-oxobutyric acid and (*rac.*)-(2*S*[,] 3*R*^{*})-4-((1*R*[,] 5*S*^{*})-6-[cyclopropyl-(3,5-dimethoxybenzyl)carbamoyl]-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-en-3-yl)-2,3-dihydroxy-4-oxobutyric acid;
- 30 2,3-dihydroxy-4-oxobutyric acid;

- 1:1-mixture of (*rac.*)-(2*R*^{*}, 3*S*^{*})-4-((1*R*^{*}, 5*S*^{*})-6-{cyclopropyl-[2-(4-methylphenyl)ethyl]carbamoyl}-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-en-3-yl)-2,3-dihydroxy-4-oxobutyric acid and (*rac.*)-(2*S*^{*}, 3*R*^{*})-4-((1*R*^{*}, 5*S*^{*})-6-{cyclopropyl-[2-(4-methylphenyl)ethyl]carbamoyl}-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo-[3.3.1]non-6-en-3-yl)-2,3-dihydroxy-4-oxobutyric acid;
- 1:1-mixture of (*rac.*)-(3*R*^{*})-5-((1*R*^{*}, 5*S*^{*})-6-[(2-chlorobenzyl)cyclopropyl-carbamoyl]-7-{4-[3-(2,3,5-trimethylphenoxy)ethyl]phenyl}-3,9-diazabicyclo-[3.3.1]non-6-en-3-yl)-3-hydroxy-5-oxopentanoic acid and (*rac.*)-(3*S*^{*})-5-((1*R*^{*}, 5*S*^{*})-6-[(2-chlorobenzyl)cyclopropyl-carbamoyl]-7-{4-[3-(2,3,5-trimethylphenoxy)ethyl]phenyl}-3,9-diazabicyclo-[3.3.1]non-6-en-3-yl)-3-hydroxy-5-oxopentanoic acid;
- 1:1-mixture of (*rac.*)-(3*R*^{*})-5-((1*R*^{*}, 5*S*^{*})-6-[(2-chlorobenzyl)ethylcarbamoyl]-7-{4-[3-(2,3,5-trimethylphenoxy)ethyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-en-3-yl)-3-hydroxy-5-oxopentanoic acid and (*rac.*)-(3*S*^{*})-5-((1*R*^{*}, 5*S*^{*})-6-[(2-chlorobenzyl)ethylcarbamoyl]-7-{4-[3-(2,3,5-trimethylphenoxy)ethyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-en-3-yl)-3-hydroxy-5-oxopentanoic acid;
- (*rac.*)-5-((1*R*^{*}, 5*S*^{*})-7-{4-[3-(2-bromo-5-fluorophenoxy)propyl]phenyl}-6-{cyclopropyl-[2-(2-hydroxyethyl)benzyl]carbamoyl}-3,9-diazabicyclo[3.3.1]non-6-en-3-yl)-5-oxo-pentanoic acid;
- (*rac.*)-5-((1*R*^{*}, 5*S*^{*})-7-{4-[3-(2-bromo-5-fluorophenoxy)propyl]phenyl}-6-{cyclopropyl-[2-(2-hydroxyethyl)benzyl]carbamoyl}-3,9-diazabicyclo[3.3.1]non-6-en-3-yl)-5-oxo-pentanoic acid methyl ester;
- (*rac.*)-(1*R*^{*}, 5*S*^{*})-7-{4-[3-(2-bromo-5-fluorophenoxy)propyl]phenyl}-3-(4-carbamoylbutyryl)-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid cyclopropyl-[2-(2-hydroxyethyl)benzyl]amide;

- (*rac.*)-(1*R*^{*,} 5*S*^{*})-3-acetyl-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid cyclopropyl-[2-(2-hydroxyethyl)-benzyl]amide;
- 5 (*rac.*)-5-((1*R*^{*,} 5*S*^{*})-6-{cyclopropyl-[2-(2-hydroxyethyl)benzyl]carbamoyl}-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-en-3-yl)-5-oxo-pentanoic acid;
- 10 (*rac.*)-5-((1*R*^{*,} 5*S*^{*})-6-{cyclopropyl-[2-(2-hydroxyethyl)benzyl]carbamoyl}-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-en-3-yl)-5-oxopentanoic acid methyl ester;
- 15 (*rac.*)-5-((1*R*^{*,} 5*S*^{*})-6-{cyclopropyl-[2-(2-hydroxyethyl)benzyl]carbamoyl}-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-en-3-yl)-2,2-dimethyl-5-oxopentanoic acid;
- 20 (*rac.*)-(1*R*^{*,} 5*S*^{*})-3-(4-carbamoylbutyryl)-7-{4-[3-(2,3,6-trifluorophenoxy)-propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid cyclopropyl-[2-(2-hydroxyethyl)benzyl]amide;
- 25 (*rac.*)-5-((1*R*^{*,} 5*S*^{*})-7-{4-[3-(2-bromo-5-fluorophenoxy)propyl]phenyl}-6-[(2-chlorobenzyl)ethylcarbamoyl]-3,9-diazabicyclo[3.3.1]non-6-en-3-yl)-5-oxo-pentanoic acid;
- 30 (*rac.*)-5-((1*R*^{*,} 5*S*^{*})-7-{4-[3-(2-bromo-5-fluorophenoxy)propyl]phenyl}-6-[cyclopropyl-(3-trifluoromethylbenzyl)carbamoyl]-3,9-diazabicyclo[3.3.1]non-6-en-3-yl)-5-oxopentanoic acid;
- 35 (*rac.*)-5-((1*R*^{*,} 5*S*^{*})-7-{4-[3-(2-bromo-5-fluorophenoxy)propyl]phenyl}-6-[cyclopropyl-(2-methylbenzyl)carbamoyl]-3,9-diazabicyclo[3.3.1]non-6-en-3-yl)-5-oxopentanoic acid;

- (*rac.*)-5-{(1*R*^{*}, 5*S*^{*})-7-[4-[3-(2-bromo-5-fluorophenoxy)propyl]phenyl]-6-[cyclopropyl-[2-(4-methoxyphenoxy)ethyl]carbamoyl]-3,9-diazabicyclo[3.3.1]-non-6-en-3-yl}-5-oxopentanoic acid;
- 5 (*rac.*)-5-{(1*R*^{*}, 5*S*^{*})-7-[4-[3-(2-bromo-5-fluorophenoxy)propyl]phenyl]-6-[cyclopropyl-[2-(3,4-dimethylphenoxy)ethyl]carbamoyl]-3,9-diazabicyclo[3.3.1]-non-6-en-3-yl}-5-oxopentanoic acid;
- 10 (*rac.*)-5-{(1*R*^{*}, 5*S*^{*})-7-[4-[3-(2-bromo-5-fluorophenoxy)propyl]phenyl]-6-[(2-(4-fluorophenyl)ethyl)cyclopropylcarbamoyl]-3,9-diazabicyclo[3.3.1]non-6-en-3-yl}-5-oxopentanoic acid;
- 15 (*rac.*)-5-{(1*R*^{*}, 5*S*^{*})-7-[4-[3-(2-bromo-5-fluorophenoxy)propyl]phenyl]-6-[(2-chlorobenzyl)ethylcarbamoyl]-3,9-diazabicyclo[3.3.1]non-6-en-3-yl}-5-oxo-pentanoic acid methyl ester;
- 20 (*rac.*)-5-{(1*R*^{*}, 5*S*^{*})-7-[4-[3-(2-bromo-5-fluorophenoxy)propyl]phenyl]-6-[cyclopropyl-(2-fluorobenzyl)carbamoyl]-3,9-diazabicyclo[3.3.1]non-6-en-3-yl}-5-oxopentanoic acid methyl ester;
- 25 (*rac.*)-5-{(1*R*^{*}, 5*S*^{*})-7-[4-[3-(2-bromo-5-fluorophenoxy)propyl]phenyl]-6-[cyclopropyl-(3-trifluoromethylbenzyl)carbamoyl]-3,9-diazabicyclo[3.3.1]non-6-en-3-yl}-5-oxopentanoic acid methyl ester;
- 30 (*rac.*)-5-{(1*R*^{*}, 5*S*^{*})-7-[4-[3-(2-bromo-5-fluorophenoxy)propyl]phenyl]-6-[cyclopropyl-[2-(4-methoxyphenoxy)ethyl]carbamoyl]-3,9-diazabicyclo[3.3.1]-non-6-en-3-yl}-5-oxopentanoic acid methyl ester;

- (*rac.*)-5-((*1R*^{*}, *5S*^{*})-7-{4-[3-(2-bromo-5-fluorophenoxy)propyl]phenyl}-6-{cyclopropyl-[2-(3-methylphenoxy)ethyl]carbamoyl}-3,9-diazabicyclo[3.3.1]non-6-en-3-yl)-5-oxopentanoic acid methyl ester;
- 5 (*rac.*)-5-((*1R*^{*}, *5S*^{*})-7-{4-[3-(2-bromo-5-fluorophenoxy)propyl]phenyl}-6-{cyclopropyl-[2-(3,4-dimethylphenoxy)ethyl]carbamoyl}-3,9-diazabicyclo[3.3.1]-non-6-en-3-yl)-5-oxopentanoic acid methyl ester;
- 10 (*rac.*)-5-[(*1R*^{*}, *5S*^{*})-7-{4-[3-(2-bromo-5-fluorophenoxy)propyl]phenyl}-6-(cyclopropylphenethylcarbamoyl)-3,9-diazabicyclo[3.3.1]non-6-en-3-yl]-5-oxo-pentanoic acid methyl ester;
- 15 (*rac.*)-5-((*1R*^{*}, *5S*^{*})-7-{4-[3-(2-bromo-5-fluorophenoxy)propyl]phenyl}-6-{[2-(2-chlorophenyl)ethyl]cyclopropylcarbamoyl}-3,9-diazabicyclo[3.3.1]non-6-en-3-yl)-5-oxopentanoic acid methyl ester;
- 20 (*rac.*)-5-((*1R*^{*}, *5S*^{*})-7-{4-[3-(2-bromo-5-fluorophenoxy)propyl]phenyl}-6-{[2-(2,3-difluorophenyl)ethyl]cyclopropylcarbamoyl}-3,9-diazabicyclo[3.3.1]non-6-en-3-yl)-5-oxopentanoic acid methyl ester;
- 25 (*rac.*)-5-((*1R*^{*}, *5S*^{*})-7-{4-[3-(2-bromo-5-fluorophenoxy)propyl]phenyl}-6-{[2-(4-fluorophenyl)ethyl]cyclopropylcarbamoyl}-3,9-diazabicyclo[3.3.1]non-6-en-3-yl)-5-oxopentanoic acid methyl ester;
- 30 (*rac.*)-5-((*1R*^{*}, *5S*^{*})-7-{4-[3-(2-bromo-5-fluorophenoxy)propyl]phenyl}-6-{[2-(4-methylphenyl)ethyl]cyclopropylcarbamoyl}-3,9-diazabicyclo[3.3.1]non-6-en-3-yl)-5-oxopentanoic acid methyl ester;

- (*rac.*)-5-{(1*R*^{*}, 5*S*^{*})-7-[4-[3-(2-bromo-5-fluorophenoxy)propyl]phenyl]-6-[(2-chlorobenzyl)ethylcarbamoyl]-3,9-diazabicyclo[3.3.1]non-6-en-3-yl}-2,2-dimethyl-5-oxopentanoic acid;
- 5 (*rac.*)-5-{(1*R*^{*}, 5*S*^{*})-7-[4-[3-(2-bromo-5-fluorophenoxy)propyl]phenyl]-6-[cyclopropyl-(3-trifluoromethylbenzyl)carbamoyl]-3,9-diazabicyclo[3.3.1]non-6-en-3-yl}-2,2-dimethyl-5-oxo-pentanoic acid;
- 10 (*rac.*)-5-{(1*R*^{*}, 5*S*^{*})-7-[4-[3-(2-bromo-5-fluorophenoxy)propyl]phenyl]-6-[cyclopropyl-[2-(4-methoxyphenoxy)ethyl]carbamoyl]-3,9-diazabicyclo[3.3.1]-non-6-en-3-yl}-2,2-dimethyl-5-oxopentanoic acid;
- 15 (*rac.*)-5-{(1*R*^{*}, 5*S*^{*})-7-[4-[3-(2-bromo-5-fluorophenoxy)propyl]phenyl]-6-[cyclopropyl-[2-(3-methylphenoxy)ethyl]carbamoyl]-3,9-diazabicyclo[3.3.1]-non-6-en-3-yl}-2,2-dimethyl-5-oxopentanoic acid;
- 20 (*rac.*)-5-{(1*R*^{*}, 5*S*^{*})-7-[4-[3-(2-bromo-5-fluorophenoxy)propyl]phenyl]-6-[cyclopropyl-[2-(3,4-dimethylphenoxy)ethyl]carbamoyl]-3,9-diazabicyclo-[3.3.1]non-6-en-3-yl}-2,2-dimethyl-5-oxopentanoic acid;
- 25 (*rac.*)-5-{(1*R*^{*}, 5*S*^{*})-7-[4-[3-(2-bromo-5-fluorophenoxy)propyl]phenyl]-6-[[2-(2-chlorophenyl)ethyl]cyclopropylcarbamoyl]-3,9-diazabicyclo[3.3.1]non-6-en-3-yl}-2,2-dimethyl-5-oxopentanoic acid;
- 30 (*rac.*)-5-{(1*R*^{*}, 5*S*^{*})-7-[4-[3-(2-bromo-5-fluorophenoxy)propyl]phenyl]-6-[[2-(4-fluorophenyl)ethyl]cyclopropylcarbamoyl]-3,9-diazabicyclo[3.3.1]non-6-en-3-yl}-2,2-dimethyl-5-oxopentanoic acid;

(*rac.*)-5-((*1R*^{*}, *5S*^{*})-7-{4-[3-(2-bromo-5-fluorophenoxy)propyl]phenyl}-6-[(2-methylphenyl)ethyl]cyclopropylcarbamoyl]-3,9-diazabicyclo[3.3.1]non-6-en-3-yl)-2,2-dimethyl-5-oxopentanoic acid;

5 1:1-mixture of (*rac.*)-(2*R*^{*}, 3*S*^{*})-4-((*1R*^{*}, *5S*^{*})-7-{4-[3-(2-bromo-5-fluorophenoxy)propyl]phenyl}-6-[(2-chlorobenzyl)ethylcarbamoyl]-3,9-diazabicyclo[3.3.1]non-6-en-3-yl}-2,3-dihydroxy-4-oxobutyric acid and (*rac.*)-(2*S*^{*}, 3*R*^{*})-4-((*1R*^{*}, *5S*^{*})-7-{4-[3-(2-bromo-5-fluorophenoxy)propyl]phenyl}-6-[(2-chlorobenzyl)ethylcarbamoyl]-3,9-diazabicyclo[3.3.1]non-6-en-3-yl}-2,3-dihydroxy-4-oxobutyric acid;
10

1:1-mixture of (*rac.*)-(2*R*^{*}, 3*S*^{*})-4-((*1R*^{*}, *5S*^{*})-7-{4-[3-(2-bromo-5-fluorophenoxy)propyl]phenyl}-6-[cyclopropyl-(2-fluorobenzyl)carbamoyl]-3,9-diazabicyclo[3.3.1]non-6-en-3-yl}-2,3-dihydroxy-4-oxobutyric acid and (*rac.*)-(2*S*^{*}, 3*R*^{*})-4-((*1R*^{*}, *5S*^{*})-7-{4-[3-(2-bromo-5-fluorophenoxy)propyl]phenyl}-6-[cyclopropyl-(2-fluorobenzyl)carbamoyl]-3,9-diazabicyclo[3.3.1]non-6-en-3-yl}-2,3-dihydroxy-4-oxobutyric acid;

1:1-mixture of (*rac.*)-(2*R*^{*}, 3*S*^{*})-4-((*1R*^{*}, *5S*^{*})-7-{4-[3-(2-bromo-5-fluorophenoxy)propyl]phenyl}-6-[cyclopropyl-(3-trifluoromethylbenzyl)carbamoyl]-3,9-diazabicyclo[3.3.1]non-6-en-3-yl}-2,3-dihydroxy-4-oxobutyric acid and (*rac.*)-(2*S*^{*}, 3*R*^{*})-4-((*1R*^{*}, *5S*^{*})-7-{4-[3-(2-bromo-5-fluorophenoxy)propyl]phenyl}-6-[cyclopropyl-(3-trifluoromethylbenzyl)carbamoyl]-3,9-diazabicyclo[3.3.1]non-6-en-3-yl}-2,3-dihydroxy-4-oxobutyric acid;
20

25 1:1-mixture of (*rac.*)-(2*R*^{*}, 3*S*^{*})-4-((*1R*^{*}, *5S*^{*})-7-{4-[3-(2-bromo-5-fluorophenoxy)propyl]phenyl}-6-[cyclopropyl-(2-methylbenzyl)carbamoyl]-3,9-diazabicyclo[3.3.1]non-6-en-3-yl}-2,3-dihydroxy-4-oxobutyric acid and (*rac.*)-(2*S*^{*}, 3*R*^{*})-4-((*1R*^{*}, *5S*^{*})-7-{4-[3-(2-bromo-5-fluorophenoxy)propyl]phenyl}-6-[cyclopropyl-(2-methylbenzyl)carbamoyl]-3,9-diazabicyclo[3.3.1]non-6-en-3-yl}-2,3-dihydroxy-4-oxobutyric acid;
30

- 1:1-mixture of (*rac.*)-(2*R*^{*,} 3*S*^{*})-4-((1*R*^{*,} 5*S*^{*})-7-{4-[3-(2-bromo-5-fluoro-phenoxy)propyl]phenyl}-6-{cyclopropyl-[2-(4-methoxyphenoxy)ethyl]-carbamoyl}-3,9-diazabicyclo[3.3.1]non-6-en-3-yl)-2,3-dihydroxy-4-oxobutyric acid and (*rac.*)-(2*S*^{*,} 3*R*^{*})-4-((1*R*^{*,} 5*S*^{*})-7-{4-[3-(2-bromo-5-fluoro-phenoxy)propyl]phenyl}-6-{cyclopropyl-[2-(4-methoxyphenoxy)ethyl]-carbamoyl}-3,9-diazabicyclo[3.3.1]non-6-en-3-yl)-2,3-dihydroxy-4-oxobutyric acid;
- 1:1-mixture of (*rac.*)-(2*R*^{*,} 3*S*^{*})-4-((1*R*^{*,} 5*S*^{*})-7-{4-[3-(2-bromo-5-fluoro-phenoxy)propyl]phenyl}-6-{cyclopropyl-[2-(3-methylphenoxy)ethyl]carbamoyl}-3,9-diazabicyclo[3.3.1]non-6-en-3-yl)-2,3-dihydroxy-4-oxobutyric acid and (*rac.*)-(2*S*^{*,} 3*R*^{*})-4-((1*R*^{*,} 5*S*^{*})-7-{4-[3-(2-bromo-5-fluorophenoxy)propyl]phenyl}-6-{cyclopropyl-[2-(3-methylphenoxy)ethyl]carbamoyl}-3,9-diazabicyclo[3.3.1]non-6-en-3-yl)-2,3-dihydroxy-4-oxobutyric acid;
- 1:1-mixture of (*rac.*)-(2*R*^{*,} 3*S*^{*})-4-((1*R*^{*,} 5*S*^{*})-7-{4-[3-(2-bromo-5-fluoro-phenoxy)propyl]phenyl}-6-{cyclopropyl-[2-(3,4-dimethylphenoxy)ethyl]-carbamoyl}-3,9-diazabicyclo[3.3.1]non-6-en-3-yl)-2,3-dihydroxy-4-oxobutyric acid and (*rac.*)-(2*S*^{*,} 3*R*^{*})-4-((1*R*^{*,} 5*S*^{*})-7-{4-[3-(2-bromo-5-fluorophenoxy)propyl]phenyl}-6-{cyclopropyl-[2-(3,4-dimethylphenoxy)ethyl]carbamoyl}-3,9-diazabicyclo[3.3.1]non-6-en-3-yl)-2,3-dihydroxy-4-oxobutyric acid;
- 1:1-mixture of (*rac.*)-(2*R*^{*,} 3*S*^{*})-4-[(1*R*^{*,} 5*S*^{*})-7-{4-[3-(2-bromo-5-fluoro-phenoxy)propyl]phenyl}-6-(cyclopropylphenethylcarbamoyl)-3,9-diazabicyclo-[3.3.1]non-6-en-3-yl]-2,3-dihydroxy-4-oxobutyric acid and (*rac.*)-(2*S*^{*,} 3*R*^{*})-4-[(1*R*^{*,} 5*S*^{*})-7-{4-[3-(2-bromo-5-fluorophenoxy)propyl]phenyl}-6-(cyclopropylphenethylcarbamoyl)-3,9-diazabicyclo-[3.3.1]non-6-en-3-yl]-2,3-dihydroxy-4-oxobutyric acid;
- 1:1-mixture of (*rac.*)-(2*R*^{*,} 3*S*^{*})-4-((1*R*^{*,} 5*S*^{*})-7-{4-[3-(2-bromo-5-fluoro-phenoxy)propyl]phenyl}-6-{[2-(2-chlorophenyl)ethyl]cyclopropylcarbamoyl}-3,9-diazabicyclo[3.3.1]non-6-en-3-yl)-2,3-dihydroxy-4-oxobutyric acid and

- (*rac.*)-(2*S*^{*}, 3*R*^{*})-4-((1*R*^{*}, 5*S*^{*})-7-{4-[3-(2-bromo-5-fluorophenoxy)propyl]-phenyl}-6-{[2-(2-chlorophenyl)ethyl]cyclopropylcarbamoyl}-3,9-diazabicyclo-[3.3.1]non-6-en-3-yl)-2,3-dihydroxy-4-oxobutyric acid;
- 5 1:1-mixture of (*rac.*)-(2*R*^{*}, 3*S*^{*})-4-((1*R*^{*}, 5*S*^{*})-7-{4-[3-(2-bromo-5-fluorophenoxy)propyl]phenyl}-6-{[2-(4-fluorophenyl)ethyl]cyclopropylcarbamoyl}-3,9-diazabicyclo[3.3.1]non-6-en-3-yl)-2,3-dihydroxy-4-oxobutyric acid and (*rac.*)-(2*S*^{*}, 3*R*^{*})-4-((1*R*^{*}, 5*S*^{*})-7-{4-[3-(2-bromo-5-fluorophenoxy)propyl]phenyl}-6-{[2-(4-fluorophenyl)ethyl]cyclopropylcarbamoyl}-3,9-diazabicyclo-[3.3.1]non-6-en-3-yl)-2,3-dihydroxy-4-oxobutyric acid;
- 10 1:1-mixture of (*rac.*)-(2*R*^{*}, 3*S*^{*})-4-((1*R*^{*}, 5*S*^{*})-7-{4-[3-(2-bromo-5-fluorophenoxy)propyl]phenyl}-6-{[2-(2-methylphenyl)ethyl]cyclopropylcarbamoyl}-3,9-diazabicyclo[3.3.1]non-6-en-3-yl)-2,3-dihydroxy-4-oxobutyric acid and (*rac.*)-(2*S*^{*}, 3*R*^{*})-4-((1*R*^{*}, 5*S*^{*})-7-{4-[3-(2-bromo-5-fluorophenoxy)propyl]phenyl}-6-{[2-(2-methylphenyl)ethyl]cyclopropylcarbamoyl}-3,9-diazabicyclo-[3.3.1]non-6-en-3-yl)-2,3-dihydroxy-4-oxobutyric acid;
- 15 1:1-mixture of (*rac.*)-(2*R*^{*}, 3*S*^{*})-4-((1*R*^{*}, 5*S*^{*})-7-{4-[3-(2-bromo-5-fluorophenoxy)propyl]phenyl}-6-{[2-(4-methylphenyl)ethyl]cyclopropylcarbamoyl}-3,9-diazabicyclo[3.3.1]non-6-en-3-yl)-2,3-dihydroxy-4-oxobutyric acid and (*rac.*)-(2*S*^{*}, 3*R*^{*})-4-((1*R*^{*}, 5*S*^{*})-7-{4-[3-(2-bromo-5-fluorophenoxy)propyl]phenyl}-6-{[2-(4-methylphenyl)ethyl]cyclopropylcarbamoyl}-3,9-diazabicyclo-[3.3.1]non-6-en-3-yl)-2,3-dihydroxy-4-oxobutyric acid;
- 20 25 (*rac.*)-(1*R*^{*}, 5*S*^{*})-7-{4-[3-(2-bromo-5-fluorophenoxy)propyl]phenyl}-3-(4-carbamoylbutyryl)-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid (2-chlorobenzyl)ethylamide;
- 30 (*rac.*)-(1*R*^{*}, 5*S*^{*})-7-{4-[3-(2-bromo-5-fluorophenoxy)propyl]phenyl}-3-(4-carbamoylbutyryl)-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid cyclopropyl-(2-fluorobenzyl)amide;

- (*rac.*)-(1*R*^{*}, 5*S*^{*})-7-{4-[3-(2-bromo-5-fluorophenoxy)propyl]phenyl}-3-(4-carbamoylbutyryl)-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid cyclopropyl-(2-methylbenzyl)amide;
- 5
- (*rac.*)-(1*R*^{*}, 5*S*^{*})-7-{4-[3-(2-bromo-5-fluorophenoxy)propyl]phenyl}-3-(4-carbamoylbutyryl)-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid cyclopropyl-[2-(4-methoxyphenoxy)ethyl]amide;
- 10 (*rac.*)-(1*R*^{*}, 5*S*^{*})-7-{4-[3-(2-bromo-5-fluorophenoxy)propyl]phenyl}-3-(4-carbamoylbutyryl)-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid cyclopropyl-[2-(3,4-dimethylphenoxy)ethyl]amide;
- 15 (*rac.*)-(1*R*^{*}, 5*S*^{*})-7-{4-[3-(2-bromo-5-fluorophenoxy)propyl]phenyl}-3-(4-carbamoylbutyryl)-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid [2-(2-chlorophenyl)ethyl]cyclopropylamide;
- 20 (*rac.*)-(1*R*^{*}, 5*S*^{*})-7-{4-[3-(2-bromo-5-fluorophenoxy)propyl]phenyl}-3-(4-carbamoylbutyryl)-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid [2-(2,3-difluorophenyl)ethyl]cyclopropylamide;
- 25 (*rac.*)-(1*R*^{*}, 5*S*^{*})-7-{4-[3-(2-bromo-5-fluorophenoxy)propyl]phenyl}-3-(4-carbamoylbutyryl)-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid [2-(4-fluorophenyl)ethyl]cyclopropylamide;
- (*rac.*)-(1*R*^{*}, 5*S*^{*})-7-{4-[3-(2-bromo-5-fluorophenoxy)propyl]phenyl}-3-(4-carbamoylbutyryl)-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid [2-(2-methylphenyl)ethyl]cyclopropylamide;
- 30 (*rac.*)-(1*R*^{*}, 5*S*^{*})-7-{4-[3-(2-bromo-5-fluorophenoxy)propyl]phenyl}-3-(4-carbamoylbutyryl)-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid [2-(4-methylphenyl)ethyl]cyclopropylamide;

(*rac.*)-(1*R**, 5*S**)-7-{4-[3-(2-bromo-5-fluorophenoxy)propyl]phenyl}-3-(4-carbamoylbutyryl)-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid cyclopropyl-(3,5-dimethoxybenzyl)amide;

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1:1-mixture of (*rac.*)-(3*R**)-5-{(*IR**, 5*S**)-7-{4-[3-(2-bromo-5-fluorophenoxy)-propyl]phenyl}-6-[(2-chlorobenzyl)ethylcarbamoyl]-3,9-diazabicyclo[3.3.1]non-6-en-3-yl}-3-hydroxy-5-oxopentanoic acid and (*rac.*)-(3*S**)-5-{(*IR**, 5*S**)-7-{4-[3-(2-bromo-5-fluorophenoxy)-propyl]phenyl}-6-[(2-chlorobenzyl)ethylcarbamoyl]-3,9-diazabicyclo[3.3.1]non-6-en-3-yl}-3-hydroxy-5-oxopentanoic acid;

1:1-mixture of (*rac.*)-(3*R**)-5-{(*IR**, 5*S**)-7-{4-[3-(2-bromo-5-fluorophenoxy)-propyl]phenyl}-6-[cyclopropyl-(2-fluorobenzyl)carbamoyl]-3,9-diazabicyclo[3.3.1]non-6-en-3-yl}-3-hydroxy-5-oxopentanoic acid and (*rac.*)-(3*S**)-5-{(*IR**, 5*S**)-7-{4-[3-(2-bromo-5-fluorophenoxy)propyl]phenyl}-6-[cyclopropyl-(2-fluorobenzyl)carbamoyl]-3,9-diazabicyclo[3.3.1]non-6-en-3-yl}-3-hydroxy-5-oxopentanoic acid;

20 1:1-mixture of (*rac.*)-(3*R**)-5-{(*IR**, 5*S**)-7-{4-[3-(2-bromo-5-fluorophenoxy)-propyl]phenyl}-6-[cyclopropyl-(3-trifluoromethylbenzyl)carbamoyl]-3,9-diazabicyclo[3.3.1]non-6-en-3-yl}-3-hydroxy-5-oxopentanoic acid and (*rac.*)-(3*S**)-5-{(*IR**, 5*S**)-7-{4-[3-(2-bromo-5-fluorophenoxy)propyl]phenyl}-6-[cyclopropyl-(3-trifluoromethylbenzyl)carbamoyl]-3,9-diazabicyclo[3.3.1]non-6-en-3-yl}-3-hydroxy-5-oxopentanoic acid;

25 1:1-mixture of (*rac.*)-(3*R**)-5-{(*IR**, 5*S**)-7-{4-[3-(2-bromo-5-fluorophenoxy)-propyl]phenyl}-6-[cyclopropyl-(2-methylbenzyl)carbamoyl]-3,9-diazabicyclo[3.3.1]non-6-en-3-yl}-3-hydroxy-5-oxopentanoic acid and (*rac.*)-(3*S**)-5-{(*IR**, 5*S**)-7-{4-[3-(2-bromo-5-fluorophenoxy)propyl]phenyl}-6-[cyclopropyl-(2-methylbenzyl)carbamoyl]-3,9-diazabicyclo[3.3.1]non-6-en-3-yl}-3-hydroxy-5-oxopentanoic acid;

- 1:1-mixture of (*rac.*)-(3*R**)-5-((1*R**, 5*S**)-7-{4-[3-(2-bromo-5-fluorophenoxy)-propyl]phenyl}-6-{cyclopropyl-[2-(4-methoxyphenoxy)ethyl]carbamoyl}-3,9-diazabicyclo[3.3.1]non-6-en-3-yl)-3-hydroxy-5-oxopentanoic acid and (*rac.*)-(3*S**)-5-((1*R**, 5*S**)-7-{4-[3-(2-bromo-5-fluorophenoxy)propyl]phenyl}-6-{cyclopropyl-[2-(4-methoxyphenoxy)ethyl]carbamoyl}-3,9-diazabicyclo[3.3.1]-non-6-en-3-yl)-3-hydroxy-5-oxopentanoic acid;
- 1:1-mixture of (*rac.*)-(3*R**)-5-((1*R**, 5*S**)-7-{4-[3-(2-bromo-5-fluorophenoxy)-propyl]phenyl}-6-{cyclopropyl-[2-(3-methylphenoxy)ethyl]carbamoyl}-3,9-diazabicyclo[3.3.1]non-6-en-3-yl)-3-hydroxy-5-oxopentanoic acid and (*rac.*)-(3*S**)-5-((1*R**, 5*S**)-7-{4-[3-(2-bromo-5-fluorophenoxy)propyl]phenyl}-6-{cyclopropyl-[2-(3-methylphenoxy)ethyl]carbamoyl}-3,9-diazabicyclo[3.3.1]non-6-en-3-yl)-3-hydroxy-5-oxopentanoic acid;
- 15 1:1-mixture of (*rac.*)-(3*R**)-5-((1*R**, 5*S**)-7-{4-[3-(2-bromo-5-fluorophenoxy)-propyl]phenyl}-6-{cyclopropyl-[2-(3,4-dimethylphenoxy)ethyl]carbamoyl}-3,9-diazabicyclo[3.3.1]non-6-en-3-yl)-3-hydroxy-5-oxopentanoic acid and (*rac.*)-(3*S**)-5-((1*R**, 5*S**)-7-{4-[3-(2-bromo-5-fluorophenoxy)propyl]phenyl}-6-{cyclopropyl-[2-(3,4-dimethylphenoxy)ethyl]carbamoyl}-3,9-diazabicyclo[3.3.1]-non-6-en-3-yl)-3-hydroxy-5-oxopentanoic acid;
- 20 1:1-mixture of (*rac.*)-(3*R**)-5-[(1*R**, 5*S**)-7-{4-[3-(2-bromo-5-fluorophenoxy)-propyl]phenyl}-6-(cyclopropylphenethylcarbamoyl)-3,9-diazabicyclo[3.3.1]non-6-en-3-yl]-3-hydroxy-5-oxopentanoic acid and (*rac.*)-(3*R**)-5-[(1*R**, 5*S**)-7-{4-[3-(2-bromo-5-fluorophenoxy)propyl]phenyl}-6-(cyclopropylphenethyl-carbamoyl)-3,9-diazabicyclo[3.3.1]non-6-en-3-yl]-3-hydroxy-5-oxopentanoic acid;
- 25 1:1-mixture of (*rac.*)-(3*R**)-5-[(1*R**, 5*S**)-7-{4-[3-(2-bromo-5-fluorophenoxy)-propyl]phenyl}-6-(cyclopropylphenethylcarbamoyl)-3,9-diazabicyclo[3.3.1]non-6-en-3-yl]-3-hydroxy-5-oxopentanoic acid and (*rac.*)-(3*R**)-5-[(1*R**, 5*S**)-7-{4-[3-(2-bromo-5-fluorophenoxy)propyl]phenyl}-6-(cyclopropylphenethyl-carbamoyl)-3,9-diazabicyclo[3.3.1]non-6-en-3-yl]-3-hydroxy-5-oxopentanoic acid;
- 30 1:1-mixture of (*rac.*)-(3*R**)-5-((1*R**, 5*S**)-7-{4-[3-(2-bromo-5-fluorophenoxy)-propyl]phenyl}-6-{{[2-(2-chlorophenyl)ethyl]cyclopropylcarbamoyl}-3,9-diaza-bicyclo[3.3.1]non-6-en-3-yl)-3-hydroxy-5-oxopentanoic acid and (*rac.*)-(3*R**)-5-

((*1R*, *5S**)-7-{4-[3-(2-bromo-5-fluorophenoxy)propyl]phenyl}-6-{[2-(2-chlorophenyl)ethyl]cyclopropylcarbamoyl}-3,9-diazabicyclo[3.3.1]non-6-en-3-yl)-3-hydroxy-5-oxopentanoic acid;

5 1:1-mixture of (*rac.*)-(3*R**)-5-((*1R*, *5S**)-7-{4-[3-(2-bromo-5-fluorophenoxy)propyl]phenyl}-6-{[2-(2,3-difluorophenyl)ethyl]cyclopropylcarbamoyl}-3,9-diazabicyclo[3.3.1]non-6-en-3-yl)-3-hydroxy-5-oxopentanoic acid and (*rac.*)-(3*R**)-5-((*1R*, *5S**)-7-{4-[3-(2-bromo-5-fluorophenoxy)propyl]phenyl}-6-{[2-(2,3-difluorophenyl)ethyl]cyclopropylcarbamoyl}-3,9-diazabicyclo[3.3.1]non-6-

10 en-3-yl)-3-hydroxy-5-oxopentanoic acid;

1:1-mixture of (*rac.*)-(3*R**)-5-((*1R*, *5S**)-7-{4-[3-(2-bromo-5-fluorophenoxy)propyl]phenyl}-6-{[2-(4-fluorophenyl)ethyl]cyclopropylcarbamoyl}-3,9-diazabicyclo[3.3.1]non-6-en-3-yl)-3-hydroxy-5-oxopentanoic acid and (*rac.*)-(3*R**)-5-((*1R*, *5S**)-7-{4-[3-(2-bromo-5-fluorophenoxy)propyl]phenyl}-6-{[2-(4-fluorophenyl)ethyl]cyclopropylcarbamoyl}-3,9-diazabicyclo[3.3.1]non-6-en-3-yl)-3-hydroxy-5-oxopentanoic acid;

1:1-mixture of (*rac.*)-(3*R**)-5-((*1R*, *5S**)-7-{4-[3-(2-bromo-5-fluorophenoxy)propyl]phenyl}-6-{[2-(2-methylphenyl)ethyl]cyclopropylcarbamoyl}-3,9-diazabicyclo[3.3.1]non-6-en-3-yl)-3-hydroxy-5-oxopentanoic acid and (*rac.*)-(3*R**)-5-((*1R*, *5S**)-7-{4-[3-(2-bromo-5-fluorophenoxy)propyl]phenyl}-6-{[2-(2-methylphenyl)ethyl]cyclopropylcarbamoyl}-3,9-diazabicyclo[3.3.1]non-6-en-3-yl)-3-hydroxy-5-oxopentanoic acid;

25 1:1-mixture of (*rac.*)-(3*R**)-5-((*1R*, *5S**)-7-{4-[3-(2-bromo-5-fluorophenoxy)propyl]phenyl}-6-{[2-(4-methylphenyl)ethyl]cyclopropylcarbamoyl}-3,9-diazabicyclo[3.3.1]non-6-en-3-yl)-3-hydroxy-5-oxopentanoic acid and (*rac.*)-(3*R**)-5-((*1R*, *5S**)-7-{4-[3-(2-bromo-5-fluorophenoxy)propyl]phenyl}-6-{[2-(4-methylphenyl)ethyl]cyclopropylcarbamoyl}-3,9-diazabicyclo[3.3.1]non-6-en-3-yl)-3-hydroxy-5-oxopentanoic acid;

- 1:1-mixture of (*rac.*)-(3*R**)-5-{(*1R**, *5S**)-7-{4-[3-(2-bromo-5-fluorophenoxy)propyl]phenyl}-6-[cyclopropyl-(3,5-dimethoxybenzyl)carbamoyl]-3,9-diaza-
bicyclo[3.3.1]non-6-en-3-yl}-3-hydroxy-5-oxopentanoic acid and (*rac.*)-(3*S**)-5-
{(*1R**, *5S**)-7-{4-[3-(2-bromo-5-fluorophenoxy)propyl]phenyl}-6-[cyclopropyl-
5 (3,5-dimethoxybenzyl)carbamoyl]-3,9-diazabicyclo[3.3.1]non-6-en-3-yl}-3-
hydroxy-5-oxopentanoic acid;
- 1:1-mixture of (*rac.*)-(3*R**)-5-((*1R**, *5S**)-6-{cyclopropyl-[2-(2-hydroxyethyl)-
benzyl]carbamoyl}-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diaza-
10 bicyclo[3.3.1]non-6-en-3-yl)-3-hydroxy-5-oxopentanoic acid and (*rac.*)-(3*S**)-5-
{(*1R**, *5S**)-6-{cyclopropyl-[2-(2-hydroxyethyl)benzyl]carbamoyl}-7-{4-[3-
(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-en-3-yl}-3-
hydroxy-5-oxopentanoic acid;
- 15 (*rac.*)-(1*R**, *5S**)-3-acetyl-7-{4-[2-(2,3,5-trimethylphenoxy)ethyl]phenyl}-3,9-
diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid (2-chlorobenzyl)cyclopropyl-
amide;
- (*rac.*)-(1*R**, *5S**)-3-acetyl-7-{4-[2-(2,3,5-trimethylphenoxy)ethyl]phenyl}-3,9-
20 diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid (2-chlorobenzyl)ethylamide;
- (*rac.*)-(1*R**, *5S**)-3-acetyl-7-{4-[2-(2,3,5-trimethylphenoxy)ethyl]phenyl}-3,9-
diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid (2-fluorobenzyl)cyclopropyl-
amide;
- 25 (*rac.*)-(1*R**, *5S**)-3-acetyl-7-{4-[2-(2,3,5-trimethylphenoxy)ethyl]phenyl}-3,9-
diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid (3-trifluoromethylbenzyl)-
cyclopropylamide;
- 30 (*rac.*)-(1*R**, *5S**)-3-acetyl-7-{4-[2-(2,3,5-trimethylphenoxy)ethyl]phenyl}-3,9-
diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid (2-methylbenzyl)cyclopropyl-
amide;

- (*rac.*)-(1*R*^{*}, 5*S*^{*})-3-acetyl-7-{4-[2-(2,3,5-trimethylphenoxy)ethyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid cyclopropyl-[2-(4-methoxyphenoxy)ethyl]amide;
- 5 (*rac.*)-5-((1*R*^{*}, 5*S*^{*})-6-(benzylcyclopropylcarbamoyl)-7-{4-[2-(2,3,5-trimethylphenoxy)ethyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-en-3-yl)-5-oxopentanoic acid;
- 10 (*rac.*)-5-((1*R*^{*}, 5*S*^{*})-6-[(2-chlorobenzyl)ethylcarbamoyl]-7-{4-[2-(2,3,5-trimethylphenoxy)ethyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-en-3-yl)-5-oxopentanoic acid;
- 15 (*rac.*)-5-((1*R*^{*}, 5*S*^{*})-6-[(2-fluorobenzyl)cyclopropylcarbamoyl]-7-{4-[2-(2,3,5-trimethylphenoxy)ethyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-en-3-yl)-5-oxo-pentanoic acid;
- 20 (*rac.*)-5-((1*R*^{*}, 5*S*^{*})-6-{cyclopropyl-[2-(4-methoxyphenoxy)ethyl]carbamoyl}-7-{4-[2-(2,3,5-trimethylphenoxy)ethyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-en-3-yl)-5-oxopentanoic acid;
- 25 (*rac.*)-5-((1*R*^{*}, 5*S*^{*})-6-{cyclopropyl-[2-(3-methoxyphenoxy)ethyl]carbamoyl}-7-{4-[2-(2,3,5-trimethylphenoxy)ethyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-en-3-yl)-5-oxopentanoic acid;
- 30 (*rac.*)-5-((1*R*^{*}, 5*S*^{*})-6-[(2-chlorobenzyl)ethylcarbamoyl]-7-{4-[2-(2,3,5-trimethylphenoxy)ethyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-en-3-yl)-5-oxopentanoic acid methyl ester;

(*rac.*)-5-((*1R*^{*}, *5S*^{*})-6-[cyclopropyl-(2-fluorobenzyl)carbamoyl]-7-{4-[2-(2,3,5-trimethylphenoxy)ethyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-en-3-yl)-5-oxo-pentanoic acid methyl ester;

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(*rac.*)-5-((*1R*^{*}, *5S*^{*})-6-[cyclopropyl-(2-methylbenzyl)carbamoyl]-7-{4-[2-(2,3,5-trimethylphenoxy)ethyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-en-3-yl)-5-oxo-pentanoic acid methyl ester;

10 (*rac.*)-5-((*1R*^{*}, *5S*^{*})-6-[(2-chlorobenzyl)cyclopropylcarbamoyl]-7-{4-[2-(2,3,5-trimethylphenoxy)ethyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-en-3-yl)-2,2-dimethyl-5-oxopentanoic acid;

15 (*rac.*)-5-((*1R*^{*}, *5S*^{*})-6-(benzylcyclopropylcarbamoyl)-7-{4-[2-(2,3,5-trimethylphenoxy)ethyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-en-3-yl)-2,2-dimethyl-5-oxopentanoic acid;

20 (*rac.*)-5-((*1R*^{*}, *5S*^{*})-6-[cyclopropyl-(2-methylbenzyl)carbamoyl]-7-{4-[2-(2,3,5-trimethylphenoxy)ethyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-en-3-yl)-2,2-dimethyl-5-oxopentanoic acid;

25 1:1-mixture of (*rac.*)-(3*R*^{*})-4-((*1R*^{*}, *5S*^{*})-6-[cyclopropyl-(2-methylbenzyl)carbamoyl]-7-{4-[2-(2,3,5-trimethylphenoxy)ethyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-en-3-yl)-2,3-dihydroxy-4-oxobutyric acid and (*rac.*)-(3*S*^{*})-4-((*1R*^{*}, *5S*^{*})-6-[cyclopropyl-(2-methylbenzyl)carbamoyl]-7-{4-[2-(2,3,5-trimethylphenoxy)ethyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-en-3-yl)-2,3-dihydroxy-4-oxobutyric acid;

30 (*rac.*)-(1*R*^{*}, *5S*^{*})-3-(4-carbamoylbutyryl)-7-{4-[2-(2,3,5-trimethylphenoxy)ethyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid benzyl-cyclopropylamide;

- (*rac.*)-(1*R*^{*}, 5*S*^{*})-3-(4-carbamoylbutyryl)-7-{4-[2-(2,3,5-trimethylphenoxy)-ethyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid (2-chlorobenzyl)ethylamide;
- 5 (*rac.*)-(1*R*^{*}, 5*S*^{*})-3-(4-carbamoylbutyryl)-7-{4-[2-(2,3,5-trimethylphenoxy)-ethyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid cyclopropyl-(2-fluorobenzyl)amide;
- 10 1:1-mixture of (1*R*, 5*S*)-7-{4-[3-(2-bromo-5-fluorophenoxy)propyl]phenyl}-3-((2*S*, 4*R*)-4-hydroxypyrrolidine-2-carbonyl)-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid cyclopropyl-(3,5-dimethoxybenzyl)amide and (1*S*, 5*R*)-7-{4-[3-(2-bromo-5-fluorophenoxy)propyl]phenyl}-3-((2*S*, 4*R*)-4-hydroxypyrrolidine-2-carbonyl)-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid cyclopropyl-(3,5-dimethoxybenzyl)amide;
- 15 (*rac.*)-5-{(1*R*^{*}, 5*S*^{*})-7-{4-[3-(2-bromo-5-fluorophenoxy)propyl]phenyl}-6-[(2-chlorobenzyl)cyclopropylcarbamoyl]-3,9-diazabicyclo[3.3.1]non-6-en-3-yl}-5-oxopentanoic acid methyl ester;
- 20 1:1-mixture of (*rac.*)-(3*R*^{*})-4-{(1*R*^{*}, 5*S*^{*})-7-{4-[3-(2-bromo-5-fluorophenoxy)propyl]phenyl}-6-[(2-chlorobenzyl)cyclopropylcarbamoyl]-3,9-diazabicyclo[3.3.1]non-6-en-3-yl}-2,3-dihydroxy-4-oxobutyric acid and (*rac.*)-(3*S*^{*})-4-{(1*R*^{*}, 5*S*^{*})-7-{4-[3-(2-bromo-5-fluorophenoxy)propyl]phenyl}-6-[(2-chlorobenzyl)cyclopropylcarbamoyl]-3,9-diazabicyclo[3.3.1]non-6-en-3-yl}-2,3-dihydroxy-4-oxobutyric acid;
- 25 (*rac.*)-5-((1*R*^{*}, 5*S*^{*})-6-(benzylcyclopropylcarbamoyl)-7-{4-[3-(2-bromo-5-fluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-en-3-yl)-5-oxopentanoic acid;
- 30 1:1-mixture of (*rac.*)-(2*R*^{*}, 3*S*^{*})-4-((1*R*^{*}, 5*S*^{*})-6-(benzylcyclopropylcarbamoyl)-7-{4-[3-(2-bromo-5-fluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-

- en-3-yl)-2,3-dihydroxy-4-oxobutyric acid and (*rac.*)-(2*S*^{*,} 3*R*^{*})-4-((1*R*^{*,} 5*S*^{*})-6-(benzylcyclopropylcarbamoyl)-7-{4-[3-(2-bromo-5-fluorophenoxy)propyl]-phenyl}-3,9-diazabicyclo[3.3.1]non-6-en-3-yl)-2,3-dihydroxy-4-oxobutyric acid;
- 5 (*rac.*)-(1*R*^{*,} 5*S*^{*})-7-{4-[3-(2-bromo-5-fluorophenoxy)propyl]phenyl}-3-(4-carbamoylbutyryl)-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid benzylcyclopropylamide;
- 10 1:1-mixture of (*rac.*)-5-{(1*R*^{*,} 5*S*^{*})-7-{4-[3-(2-bromo-5-fluorophenoxy)propyl]-phenyl}-6-[(2*R*^{*})-2-hydroxy-2-phenylethyl)methylcarbamoyl]-3,9-diazabicyclo-[3.3.1]non-6-en-3-yl}-5-oxopentanoic acid and (*rac.*)-5-{(1*R*^{*,} 5*S*^{*})-7-{4-[3-(2-bromo-5-fluorophenoxy)propyl]phenyl}-6-[(2*S*^{*})-2-hydroxy-2-phenylethyl)-methylcarbamoyl]-3,9-diazabicyclo-[3.3.1]non-6-en-3-yl}-5-oxopentanoic acid;
- 15 1:1-mixture of (*rac.*)-5-{(1*R*^{*,} 5*S*^{*})-7-{4-[3-(2-bromo-5-fluorophenoxy)propyl]-phenyl}-6-[(2*R*^{*})-2-hydroxy-2-phenylethyl)methylcarbamoyl]-3,9-diazabicyclo-[3.3.1]non-6-en-3-yl}-5-oxopentanoic acid methyl ester and (*rac.*)-5-{(1*R*^{*,} 5*S*^{*})-7-{4-[3-(2-bromo-5-fluorophenoxy)propyl]phenyl}-6-[(2*S*^{*})-2-hydroxy-2-phenylethyl)methylcarbamoyl]-3,9-diazabicyclo-[3.3.1]non-6-en-3-yl}-5-oxopentanoic acid methyl ester;
- 20 1:1-mixture of (*rac.*)-5-{(1*R*^{*,} 5*S*^{*})-7-{4-[3-(2-bromo-5-fluorophenoxy)propyl]-phenyl}-6-[(2*R*^{*})-2-hydroxy-2-phenylethyl)methylcarbamoyl]-3,9-diazabicyclo-[3.3.1]non-6-en-3-yl}-2,2-dimethyl-5-oxopentanoic acid and (*rac.*)-5-{(1*R*^{*,} 5*S*^{*})-7-{4-[3-(2-bromo-5-fluorophenoxy)propyl]phenyl}-6-[(2*S*^{*})-2-hydroxy-2-phenylethyl)methylcarbamoyl]-3,9-diazabicyclo-[3.3.1]non-6-en-3-yl}-2,2-dimethyl-5-oxopentanoic acid;
- 25 1:1-mixture of (*rac.*)-(1*R*^{*,} 5*S*^{*})-7-{4-[3-(2-bromo-5-fluorophenoxy)propyl]-phenyl}-3-(4-carbamoylbutyryl)-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid ((2*R*^{*})-2-hydroxy-2-phenylethyl)methylamide and (*rac.*)-(1*R*^{*,} 5*S*^{*})-7-{4-[3-(2-bromo-5-fluorophenoxy)propyl]-phenyl}-3-(4-carbamoylbutyryl)-3,9-

diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid ((2*R**)-2-hydroxy-2-phenylethyl)-methylamide;

1:1-mixture of (*rac.*)-(3*R**)-5-((*IR**, *5S**)-7-{4-[3-(2-bromo-5-fluorophenoxy)propyl]phenyl}-6-[(2-chlorobenzyl)cyclopropylcarbamoyl]-3,9-diazabicyclo-[3.3.1]non-6-en-3-yl}-3-hydroxy-5-oxopentanoic acid and (*rac.*)-(3*S**)-5-((*IR**, *5S**)-7-{4-[3-(2-bromo-5-fluorophenoxy)propyl]phenyl}-6-[(2-chlorobenzyl)cyclopropylcarbamoyl]-3,9-diazabicyclo-[3.3.1]non-6-en-3-yl}-3-hydroxy-5-oxopentanoic acid;

10

1:1-mixture of (*rac.*)-(3*R**)-5-((*IR**, *5S**)-6-(benzylcyclopropylcarbamoyl)-7-{4-[3-(2-bromo-5-fluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-en-3-yl}-3-hydroxy-5-oxopentanoic acid and (*rac.*)-(3*S**)-5-((*IR**, *5S**)-6-(benzylcyclopropylcarbamoyl)-7-{4-[3-(2-bromo-5-fluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-en-3-yl)-3-hydroxy-5-oxopentanoic acid;

15

(*rac.*)-(1*R**, *5S**)-7-{4-[2-(2-chloro-4,5-dimethylphenoxy)ethoxy]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid [2-(2-chlorophenyl)ethyl]-cyclopropylamide;

20

(*rac.*)-(1*R**, *5S**)-7-{4-[2-(2-chloro-4,5-dimethylphenoxy)ethoxy]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid cyclopropyl-(3,5-dimethoxybenzyl)amide;

25

(*rac.*)-(1*R**, *5S**)-7-{4-[2-(2-chloro-4,5-dimethylphenoxy)ethoxy]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid cyclopropyl-(*p*-toluyl-ethyl)amide;

30

(*rac.*)-(1*R**, *5S**)-3-acetyl-7-{4-[2-(4-bromophenoxy)ethoxy]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid (2-chlorobenzyl)cyclopropylamide;

(*rac.*)-(1*R*^{*,} 5*S*^{*})-3-acetyl-7-{4-[2-(4-bromophenoxy)ethoxy]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid cyclopropyl-(3,5-dimethoxybenzyl)-amide;

5 (*rac.*)-5-{(*1R*^{*,} 5*S*^{*})-7-{4-[2-(4-bromophenoxy)ethoxy]phenyl}-6-[(2-chlorobenzyl)cyclopropylcarbamoyl]-3,9-diazabicyclo[3.3.1]non-6-en-3-yl}-5-oxo-pentanoic acid;

10 (*rac.*)-5-((*1R*^{*,} 5*S*^{*})-6-(benzylcyclopropylcarbamoyl)-7-{4-[2-(4-bromophenoxy)ethoxy]phenyl}-3,9-diazabicyclo[3.3.1]non-6-en-3-yl)-5-oxopentanoic acid;

15 (*rac.*)-5-{(*1R*^{*,} 5*S*^{*})-7-{4-[2-(4-bromophenoxy)ethoxy]phenyl}-6-[cyclopropyl-(2-fluorobenzyl)carbamoyl]-3,9-diazabicyclo[3.3.1]non-6-en-3-yl}-5-oxo-pentanoic acid;

15 (*rac.*)-5-{(*1R*^{*,} 5*S*^{*})-7-{4-[2-(4-bromophenoxy)ethoxy]phenyl}-6-[cyclopropyl-(3-trifluoromethylbenzyl)carbamoyl]-3,9-diazabicyclo[3.3.1]non-6-en-3-yl}-5-oxopentanoic acid;

20 (*rac.*)-5-{(*1R*^{*,} 5*S*^{*})-7-{4-[2-(4-bromophenoxy)ethoxy]phenyl}-6-[cyclopropyl-(2-methylbenzyl)carbamoyl]-3,9-diazabicyclo[3.3.1]non-6-en-3-yl}-5-oxo-pentanoic acid;

25 (*rac.*)-5-{(*1R*^{*,} 5*S*^{*})-7-{4-[2-(4-bromophenoxy)ethoxy]phenyl}-6-[cyclopropyl-[2-(3,4-dimethylphenoxy)ethyl]carbamoyl]-3,9-diazabicyclo[3.3.1]non-6-en-3-yl}-5-oxopentanoic acid;

30 (*rac.*)-5-{(*1R*^{*,} 5*S*^{*})-7-{4-[2-(4-bromophenoxy)ethoxy]phenyl}-6-{{[2-(2-chlorophenyl)ethyl]cyclopropylcarbamoyl}-3,9-diazabicyclo[3.3.1]non-6-en-3-yl}-5-oxopentanoic acid;

- (*rac.*)-5-{(1*R*^{*}, 5*S*^{*})-7-[4-[2-(4-bromophenoxy)ethoxy]phenyl]-6-[cyclopropyl-[2-(4-fluorophenyl)ethyl]carbamoyl]-3,9-diazabicyclo[3.3.1]non-6-en-3-yl}-5-oxopentanoic acid;
- 5 (*rac.*)-5-{(1*R*^{*}, 5*S*^{*})-7-[4-[2-(4-bromophenoxy)ethoxy]phenyl]-6-[cyclopropyl-[2-(2-methylphenyl)ethyl]carbamoyl]-3,9-diazabicyclo[3.3.1]non-6-en-3-yl}-5-oxopentanoic acid;
- 10 (*rac.*)-5-{(1*R*^{*}, 5*S*^{*})-7-[4-[2-(4-bromophenoxy)ethoxy]phenyl]-6-[cyclopropyl-[2-(4-methylphenyl)ethyl]carbamoyl]-3,9-diazabicyclo[3.3.1]non-6-en-3-yl}-5-oxopentanoic acid;
- 15 (*rac.*)-5-{(1*R*^{*}, 5*S*^{*})-7-[4-[2-(4-bromophenoxy)ethoxy]phenyl]-6-[cyclopropyl-(3,5-dimethoxybenzyl)carbamoyl]-3,9-diazabicyclo[3.3.1]non-6-en-3-yl}-5-oxo-pentanoic acid;
- 20 (*rac.*)-5-{(1*R*^{*}, 5*S*^{*})-7-[4-[2-(4-bromophenoxy)ethoxy]phenyl]-6-[cyclopropyl-(2-fluorobenzyl)carbamoyl]-3,9-diazabicyclo[3.3.1]non-6-en-3-yl}-2,2-dimethyl-5-oxopentanoic acid;
- 25 (*rac.*)-5-{(1*R*^{*}, 5*S*^{*})-7-[4-[2-(4-bromophenoxy)ethoxy]phenyl]-6-[[2-(2-chlorophenyl)ethyl]cyclopropylcarbamoyl]-3,9-diazabicyclo[3.3.1]non-6-en-3-yl}-2,2-dimethyl-5-oxopentanoic acid;
- 30 (*rac.*)-5-{(1*R*^{*}, 5*S*^{*})-7-[4-[2-(4-bromophenoxy)ethoxy]phenyl]-6-[cyclopropyl-[2-(2,3-difluorophenyl)ethyl]carbamoyl]-3,9-diazabicyclo[3.3.1]non-6-en-3-yl}-2,2-dimethyl-5-oxopentanoic acid;
- 35 (*rac.*)-5-{(1*R*^{*}, 5*S*^{*})-7-[4-[2-(4-bromophenoxy)ethoxy]phenyl]-6-[cyclopropyl-[2-(2-methylphenyl)ethyl]carbamoyl]-3,9-diazabicyclo[3.3.1]non-6-en-3-yl}-2,2-dimethyl-5-oxopentanoic acid;

- (*rac.*)-5-((*1R*^{*}, *5S*^{*})-7-{4-[2-(4-bromophenoxy)ethoxy]phenyl}-6-{cyclopropyl-[2-(4-methylphenyl)ethyl]carbamoyl}-3,9-diazabicyclo[3.3.1]non-6-en-3-yl)-2,2-dimethyl-5-oxopentanoic acid;
- 5 (*rac.*)-5-((*1R*^{*}, *5S*^{*})-7-{4-[2-(4-bromophenoxy)ethoxy]phenyl}-6-[cyclopropyl-(3,5-dimethoxybenzyl)carbamoyl]-3,9-diazabicyclo[3.3.1]non-6-en-3-yl)-2,2-dimethyl-5-oxopentanoic acid;
- 10 (*rac.*)-(1*R*^{*}, *5S*^{*})-7-{4-[2-(4-bromophenoxy)ethoxy]phenyl}-3-(4-carbamoylbutyryl)-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid benzylcyclopropylamide;
- 15 (*rac.*)-(1*R*^{*}, *5S*^{*})-7-{4-[2-(4-bromophenoxy)ethoxy]phenyl}-3-(4-carbamoylbutyryl)-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid (2-chlorobenzyl)-ethylamide;
- 20 (*rac.*)-(1*R*^{*}, *5S*^{*})-7-{4-[2-(4-bromophenoxy)ethoxy]phenyl}-3-(4-carbamoylbutyryl)-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid cyclopropyl-(2-fluorobenzyl)amide;
- 25 (*rac.*)-(1*R*^{*}, *5S*^{*})-7-{4-[2-(4-bromophenoxy)ethoxy]phenyl}-3-(4-carbamoylbutyryl)-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid cyclopropyl-[2-(4-methoxyphenoxy)ethyl]amide;
- 30 (*rac.*)-(1*R*^{*}, *5S*^{*})-7-{4-[2-(4-bromophenoxy)ethoxy]phenyl}-3-(4-carbamoylbutyryl)-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid [2-(2-chlorophenyl)-ethyl]cyclopropylamide;
- 35 (*rac.*)-(1*R*^{*}, *5S*^{*})-7-{4-[2-(4-bromophenoxy)ethoxy]phenyl}-3-(4-carbamoylbutyryl)-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid cyclopropyl-(3,5-dimethoxybenzyl)amide;

- (*rac.*)-(1*R*^{*}, 5*S*^{*})-7-{4-[2-(4-bromophenoxy)ethoxy]phenyl}-3-(4-carbamoylbutyryl)-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid cyclopropyl-[2-(2-hydroxyethyl)benzyl]amide;
- 5 (*rac.*)-(1*R*^{*}, 5*S*^{*})-7-{4-[2-(2-chloro-4,5-dimethylphenoxy)ethoxy]phenyl}-3-formyl-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid (2-chlorobenzyl)-ethylamide;
- 10 (*rac.*)-(1*R*^{*}, 5*S*^{*})-7-{4-[2-(2-chloro-4,5-dimethylphenoxy)ethoxy]phenyl}-3-formyl-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid cyclopropyl-(2-*o*-tolylethyl)amide;
- 15 (*rac.*)-(1*R*^{*}, 5*S*^{*})-7-{4-[2-(2-chloro-4,5-dimethylphenoxy)ethoxy]phenyl}-3-formyl-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid cyclopropyl-(3,5-dimethoxybenzyl)amide;
- 20 (*rac.*)-(1*R*^{*}, 5*S*^{*})-3-acetyl-7-{4-[2-(2-chloro-4,5-dimethylphenoxy)ethoxy]-phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid (2-chlorobenzyl)-ethylamide;
- 25 (*rac.*)-(1*R*^{*}, 5*S*^{*})-3-acetyl-7-{4-[2-(2-chloro-4,5-dimethylphenoxy)ethoxy]-phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid cyclopropyl-(2-fluorobenzyl)amide;
- 30 (*rac.*)-(1*R*^{*}, 5*S*^{*})-3-acetyl-7-{4-[2-(2-chloro-4,5-dimethylphenoxy)ethoxy]-phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid cyclopropyl-(2-methylbenzyl)amide;
- 35 (*rac.*)-(1*R*^{*}, 5*S*^{*})-3-acetyl-7-{4-[2-(2-chloro-4,5-dimethylphenoxy)ethoxy]-phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid cyclopropyl-(3,5-dimethoxybenzyl)amide;

- (*rac.*)-(1*R*^{*}, 5*S*^{*})-3-acetyl-7-{4-[2-(2-chloro-4,5-dimethylphenoxy)ethoxy]-phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid cyclopropyl-(2-*p*-tolylethyl)amide;
- 5 (*rac.*)-5-((1*R*^{*}, 5*S*^{*})-6-[(2-chlorobenzyl)ethylcarbamoyl]-7-{4-[2-(2-chloro-4,5-dimethylphenoxy)ethoxy]phenyl}-3,9-diazabicyclo[3.3.1]non-6-en-3-yl)-5-oxopentanoic acid;
- 10 (*rac.*)-5-((1*R*^{*}, 5*S*^{*})-7-{4-[2-(2-chloro-4,5-dimethylphenoxy)ethoxy]phenyl}-6-[cyclopropyl-(2-fluorobenzyl)carbamoyl]-3,9-diazabicyclo[3.3.1]non-6-en-3-yl)-5-oxopentanoic acid;
- 15 (*rac.*)-5-((1*R*^{*}, 5*S*^{*})-7-{4-[2-(2-chloro-4,5-dimethylphenoxy)ethoxy]phenyl}-6-[cyclopropyl-(3-trifluoromethylbenzyl)carbamoyl]-3,9-diazabicyclo[3.3.1]non-6-en-3-yl)-5-oxopentanoic acid;
- 20 (*rac.*)-5-((1*R*^{*}, 5*S*^{*})-7-{4-[2-(2-chloro-4,5-dimethylphenoxy)ethoxy]phenyl}-6-{cyclopropyl-[2-(4-methoxyphenoxy)ethyl]carbamoyl}-3,9-diazabicyclo[3.3.1]-non-6-en-3-yl)-5-oxopentanoic acid;
- 25 (*rac.*)-5-((1*R*^{*}, 5*S*^{*})-7-{4-[2-(2-chloro-4,5-dimethylphenoxy)ethoxy]phenyl}-6-{cyclopropyl-[2-(3-methoxyphenoxy)ethyl]carbamoyl}-3,9-diazabicyclo[3.3.1]-non-6-en-3-yl)-5-oxopentanoic acid;
- 30 (*rac.*)-5-((1*R*^{*}, 5*S*^{*})-7-{4-[2-(2-chloro-4,5-dimethylphenoxy)ethoxy]phenyl}-6-(cyclopropylphenethylcarbamoyl)-3,9-diazabicyclo[3.3.1]non-6-en-3-yl)-5-oxopentanoic acid;

- (*rac.*)-5-((*1R*^{*}, *5S*^{*})-7-{4-[2-(2-chloro-4,5-dimethylphenoxy)ethoxy]phenyl}-6-
{[2-(2-chlorophenyl)ethyl]cyclopropyl-carbamoyl}-3,9-diazabicyclo[3.3.1]non-6-
en-3-yl)-5-oxopentanoic acid;
- 5 (*rac.*)-5-((*1R*^{*}, *5S*^{*})-7-{4-[2-(2-chloro-4,5-dimethylphenoxy)ethoxy]phenyl}-6-
{cyclopropyl-[2-(2,3-difluorophenyl)ethyl]carbamoyl}-3,9-diazabicyclo[3.3.1]-
non-6-en-3-yl)-5-oxopentanoic acid;
- 10 (*rac.*)-5-((*1R*^{*}, *5S*^{*})-7-{4-[2-(2-chloro-4,5-dimethylphenoxy)ethoxy]phenyl}-6-
{cyclopropyl-[2-(4-fluorophenyl)ethyl]carbamoyl}-3,9-diazabicyclo[3.3.1]-non-6-
en-3-yl)-5-oxopentanoic acid;
- 15 (*rac.*)-5-((*1R*^{*}, *5S*^{*})-7-{4-[2-(2-chloro-4,5-dimethylphenoxy)ethoxy]phenyl}-6-
{cyclopropyl-[2-(2-methylphenyl)ethyl]carbamoyl}-3,9-diazabicyclo[3.3.1]-non-
6-en-3-yl)-5-oxopentanoic acid;
- 20 (*rac.*)-5-((*1R*^{*}, *5S*^{*})-7-{4-[2-(2-chloro-4,5-dimethylphenoxy)ethoxy]phenyl}-6-
{cyclopropyl-[2-(4-methylphenyl)ethyl]carbamoyl}-3,9-diazabicyclo[3.3.1]-non-
6-en-3-yl)-5-oxopentanoic acid;
- 25 (*rac.*)-5-((*1R*^{*}, *5S*^{*})-6-[(2-chlorobenzyl)cyclopropylcarbamoyl]-7-{4-[2-(2-
chloro-4,5-dimethylphenoxy)ethoxy]phenyl}-3,9-diazabicyclo[3.3.1]non-6-en-3-
yl)-5-oxopentanoic acid methyl ester;
- 30 (*rac.*)-5-((*1R*^{*}, *5S*^{*})-6-[(2-chlorobenzyl)ethylcarbamoyl]-7-{4-[2-(2-chloro-4,5-
dimethylphenoxy)ethoxy]phenyl}-3,9-diazabicyclo[3.3.1]non-6-en-3-yl)-5-
oxopentanoic acid methyl ester;
- 35 (*rac.*)-5-((*1R*^{*}, *5S*^{*})-7-{4-[2-(2-chloro-4,5-dimethylphenoxy)ethoxy]phenyl}-6-
[cyclopropyl-(2-fluorobenzyl)carbamoyl]-3,9-diazabicyclo[3.3.1]non-6-en-3-yl)-
5-oxopentanoic acid methyl ester;

- (*rac.*)-5-{(*1R*^{*}, *5S*^{*})-7-{4-[2-(2-chloro-4,5-dimethylphenoxy)ethoxy]phenyl}-6-[cyclopropyl-(2-methylbenzyl)carbamoyl]-3,9-diazabicyclo[3.3.1]non-6-en-3-yl}-5-oxopentanoic acid methyl ester;
- 5 (*rac.*)-5-{(*1R*^{*}, *5S*^{*})-7-{4-[2-(2-chloro-4,5-dimethylphenoxy)ethoxy]phenyl}-6-[cyclopropyl-(3-trifluoromethylbenzyl)carbamoyl]-3,9-diazabicyclo[3.3.1]non-6-en-3-yl}-5-oxopentanoic acid methyl ester;
- 10 (*rac.*)-5-[(*1R*^{*}, *5S*^{*})-7-{4-[2-(2-chloro-4,5-dimethylphenoxy)ethoxy]phenyl}-6-(cyclopropylphenethylcarbamoyl)-3,9-diazabicyclo[3.3.1]non-6-en-3-yl]-5-oxo-pentanoic acid methyl ester;
- 15 (*rac.*)-5-((*1R*^{*}, *5S*^{*})-7-{4-[2-(2-chloro-4,5-dimethylphenoxy)ethoxy]phenyl}-6-[[2-(2-chlorophenyl)ethyl]cyclopropylcarbamoyl]-3,9-diazabicyclo[3.3.1]non-6-en-3-yl)-5-oxopentanoic acid methyl ester;
- 20 (*rac.*)-5-((*1R*^{*}, *5S*^{*})-7-{4-[2-(2-chloro-4,5-dimethylphenoxy)ethoxy]phenyl}-6-[[2-(2,3-difluorophenyl)ethyl]cyclopropylcarbamoyl]-3,9-diazabicyclo[3.3.1]non-6-en-3-yl)-5-oxopentanoic acid methyl ester;
- 25 (*rac.*)-5-((*1R*^{*}, *5S*^{*})-7-{4-[2-(2-chloro-4,5-dimethylphenoxy)ethoxy]phenyl}-6-[[2-(4-fluorophenyl)ethyl]cyclopropylcarbamoyl]-3,9-diazabicyclo[3.3.1]non-6-en-3-yl)-5-oxopentanoic acid methyl ester;
- 30 (*rac.*)-5-((*1R*^{*}, *5S*^{*})-7-{4-[2-(2-chloro-4,5-dimethylphenoxy)ethoxy]phenyl}-6-[[2-(4-methylphenyl)ethyl]cyclopropylcarbamoyl]-3,9-diazabicyclo[3.3.1]non-6-en-3-yl)-5-oxopentanoic acid methyl ester;

- (*rac.*)-5-{(*1R*^{*}, *5S*^{*})-7-{4-[2-(2-chloro-4,5-dimethylphenoxy)ethoxy]phenyl}-6-[cyclopropyl-(3,5-dimethoxybenzyl)carbamoyl]-3,9-diazabicyclo[3.3.1]non-6-en-3-yl}-5-oxopentanoic acid methyl ester;
- 5 (*rac.*)-5-{(*1R*^{*}, *5S*^{*})-6-(benzylcyclopropylcarbamoyl)-7-{4-[2-(2-chloro-4,5-dimethylphenoxy)ethoxy]phenyl}-3,9-diazabicyclo[3.3.1]non-6-en-3-yl)-2,2-dimethyl-5-oxopentanoic acid;
- 10 (*rac.*)-5-{(*1R*^{*}, *5S*^{*})-7-{4-[2-(2-chloro-4,5-dimethylphenoxy)ethoxy]phenyl}-6-[cyclopropyl-(2-methylbenzyl)carbamoyl]-3,9-diazabicyclo[3.3.1]non-6-en-3-yl}-2,2-dimethyl-5-oxopentanoic acid;
- 15 (*rac.*)-5-{(*1R*^{*}, *5S*^{*})-7-{4-[2-(2-chloro-4,5-dimethylphenoxy)ethoxy]phenyl}-6-[cyclopropyl-[2-(3,4-dimethylphenoxy)ethyl]carbamoyl]-3,9-diazabicyclo[3.3.1]non-6-en-3-yl)-2,2-dimethyl-5-oxopentanoic acid;
- 20 (*rac.*)-5-{(*1R*^{*}, *5S*^{*})-7-{4-[2-(2-chloro-4,5-dimethylphenoxy)ethoxy]phenyl}-6-(cyclopropylphenethylcarbamoyl)-3,9-diazabicyclo[3.3.1]non-6-en-3-yl)-2,2-dimethyl-5-oxopentanoic acid;
- 25 (*rac.*)-5-{(*1R*^{*}, *5S*^{*})-7-{4-[2-(2-chloro-4,5-dimethylphenoxy)ethoxy]phenyl}-6-[[2-(2-chlorophenyl)ethyl]cyclopropylcarbamoyl]-3,9-diazabicyclo[3.3.1]non-6-en-3-yl)-2,2-dimethyl-5-oxopentanoic acid;
- 30 (*rac.*)-5-{(*1R*^{*}, *5S*^{*})-7-{4-[2-(2-chloro-4,5-dimethylphenoxy)ethoxy]phenyl}-6-[[2-(2,3-difluorophenyl)ethyl]cyclopropylcarbamoyl]-3,9-diazabicyclo[3.3.1]non-6-en-3-yl)-2,2-dimethyl-5-oxopentanoic acid;
- 35 (*rac.*)-5-{(*1R*^{*}, *5S*^{*})-7-{4-[2-(2-chloro-4,5-dimethylphenoxy)ethoxy]phenyl}-6-[[2-(4-fluorophenyl)ethyl]cyclopropylcarbamoyl]-3,9-diazabicyclo[3.3.1]non-6-en-3-yl)-2,2-dimethyl-5-oxopentanoic acid;

- (*rac.*)-5-((*1R*^{*}, *5S*^{*})-7-{4-[2-(2-chloro-4,5-dimethylphenoxy)ethoxy]phenyl}-6-
{[2-(2-methylphenyl)ethyl]cyclopropylcarbamoyl}-3,9-diazabicyclo[3.3.1]non-6-
en-3-yl)-2,2-dimethyl-5-oxopentanoic acid;
- 5 (*rac.*)-5-((*1R*^{*}, *5S*^{*})-7-{4-[2-(2-chloro-4,5-dimethylphenoxy)ethoxy]phenyl}-6-
{[2-(4-methylphenyl)ethyl]cyclopropylcarbamoyl}-3,9-diazabicyclo[3.3.1]non-6-
en-3-yl)-2,2-dimethyl-5-oxopentanoic acid;
- 10 (*rac.*)-5-((*1R*^{*}, *5S*^{*})-7-{4-[2-(2-chloro-4,5-dimethylphenoxy)ethoxy]phenyl}-6-
[cyclopropyl-(3,5-dimethoxybenzyl)carbamoyl]-3,9-diazabicyclo[3.3.1]non-6-en-
3-yl)-2,2-dimethyl-5-oxopentanoic acid;
- 15 (*rac.*)-(*1R*^{*}, *5S*^{*})-3-(4-carbamoylbutyryl)-7-{4-[2-(2-chloro-4,5-dimethyl-
phenoxy)ethoxy]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid (2-
chlorobenzyl)cyclopropylamide;
- 20 (*rac.*)-(*1R*^{*}, *5S*^{*})-3-(4-carbamoylbutyryl)-7-{4-[2-(2-chloro-4,5-dimethyl-
phenoxy)ethoxy]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid
benzylcyclopropylamide;
- 25 (*rac.*)-(*1R*^{*}, *5S*^{*})-3-(4-carbamoylbutyryl)-7-{4-[2-(2-chloro-4,5-dimethyl-
phenoxy)ethoxy]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid (2-
chlorobenzyl)ethylamide;
- 30 (*rac.*)-(*1R*^{*}, *5S*^{*})-3-(4-carbamoylbutyryl)-7-{4-[2-(2-chloro-4,5-dimethyl-
phenoxy)ethoxy]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid
cyclopropyl-(2-fluorobenzyl)amide;
- 30 (*rac.*)-(*1R*^{*}, *5S*^{*})-3-(4-carbamoylbutyryl)-7-{4-[2-(2-chloro-4,5-dimethyl-
phenoxy)ethoxy]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid
cyclopropyl-(3-trifluoromethylbenzyl)amide;

- (*rac.*)-(1*R*^{*}, 5*S*^{*})-3-(4-carbamoylbutyryl)-7-{4-[2-(2-chloro-4,5-dimethylphenoxy)ethoxy]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid cyclopropyl-(2-methylbenzyl)amide;
- 5 (*rac.*)-(1*R*^{*}, 5*S*^{*})-3-(4-carbamoylbutyryl)-7-{4-[2-(2-chloro-4,5-dimethylphenoxy)ethoxy]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid cyclopropyl-[2-(4-methoxyphenoxy)ethyl]amide;
- 10 (*rac.*)-(1*R*^{*}, 5*S*^{*})-3-(4-carbamoylbutyryl)-7-{4-[2-(2-chloro-4,5-dimethylphenoxy)ethoxy]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid cyclopropylphenethylamide;
- 15 (*rac.*)-(1*R*^{*}, 5*S*^{*})-3-(4-carbamoylbutyryl)-7-{4-[2-(2-chloro-4,5-dimethylphenoxy)ethoxy]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid [2-(2-chlorophenyl)ethyl]cyclopropylamide;
- 20 (*rac.*)-(1*R*^{*}, 5*S*^{*})-3-(4-carbamoylbutyryl)-7-{4-[2-(2-chloro-4,5-dimethylphenoxy)ethoxy]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid cyclopropyl-[2-(2,3-difluorophenyl)ethyl]amide;
- 25 (*rac.*)-(1*R*^{*}, 5*S*^{*})-3-(4-carbamoylbutyryl)-7-{4-[2-(2-chloro-4,5-dimethylphenoxy)ethoxy]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid cyclopropyl-[2-(4-fluorophenyl)ethyl]amide;
- 30 (*rac.*)-(1*R*^{*}, 5*S*^{*})-3-(4-carbamoylbutyryl)-7-{4-[2-(2-chloro-4,5-dimethylphenoxy)ethoxy]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid cyclopropyl-(2-*p*-tolylethyl)amide;

(*rac.*)-(1*R*^{*}, 5*S*^{*})-3-(4-carbamoylbutyryl)-7-{4-[2-(2-chloro-4,5-dimethylphenoxy)ethoxy]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid cyclopropyl-(3,5-dimethoxybenzyl)amide;

5 1:1-mixture of (*rac.*)-(3*R*^{*})-5-((1*R*^{*}, 5*S*^{*})-6-[(2-chlorobenzyl)cyclopropylcarbamoyl]-7-{4-[2-(2-chloro-4,5-dimethylphenoxy)ethoxy]phenyl}-3,9-diazabicyclo[3.3.1]non-6-en-3-yl)-3-hydroxy-5-oxopentanoic acid and (*rac.*)-(3*S*^{*})-5-((1*R*^{*}, 5*S*^{*})-6-[(2-chlorobenzyl)cyclopropylcarbamoyl]-7-{4-[2-(2-chloro-4,5-dimethylphenoxy)ethoxy]phenyl}-3,9-diazabicyclo[3.3.1]non-6-en-3-yl)-3-hydroxy-5-oxopentanoic acid;

10 1:1-mixture of (*rac.*)-(3*R*^{*})-5-((1*R*^{*}, 5*S*^{*})-6-[(2-chlorobenzyl)ethylcarbamoyl]-7-{4-[2-(2-chloro-4,5-dimethylphenoxy)ethoxy]phenyl}-3,9-diazabicyclo[3.3.1]non-6-en-3-yl)-3-hydroxy-5-oxopentanoic acid and (*rac.*)-(3*S*^{*})-5-((1*R*^{*}, 5*S*^{*})-6-[(2-chlorobenzyl)ethylcarbamoyl]-7-{4-[2-(2-chloro-4,5-dimethylphenoxy)ethoxy]phenyl}-3,9-diazabicyclo[3.3.1]non-6-en-3-yl)-3-hydroxy-5-oxopentanoic acid;

15 1:1-mixture of (*rac.*)-(3*R*^{*})-5-((1*R*^{*}, 5*S*^{*})-7-{4-[2-(2-chloro-4,5-dimethylphenoxy)ethoxy]phenyl}-6-[cyclopropyl-(2-fluorobenzyl)carbamoyl]-3,9-diazabicyclo[3.3.1]non-6-en-3-yl)-3-hydroxy-5-oxopentanoic acid and (*rac.*)-(3*S*^{*})-5-((1*R*^{*}, 5*S*^{*})-7-{4-[2-(2-chloro-4,5-dimethylphenoxy)ethoxy]phenyl}-6-[cyclopropyl-(2-fluorobenzyl)carbamoyl]-3,9-diazabicyclo[3.3.1]non-6-en-3-yl)-3-hydroxy-5-oxopentanoic acid;

20 1:1-mixture of (*rac.*)-(3*R*^{*})-5-((1*R*^{*}, 5*S*^{*})-7-{4-[2-(2-chloro-4,5-dimethylphenoxy)ethoxy]phenyl}-6-[cyclopropyl-(3-trifluoromethylbenzyl)carbamoyl]-3,9-diazabicyclo[3.3.1]non-6-en-3-yl)-3-hydroxy-5-oxopentanoic acid and (*rac.*)-(3*S*^{*})-5-((1*R*^{*}, 5*S*^{*})-7-{4-[2-(2-chloro-4,5-dimethylphenoxy)ethoxy]phenyl}-6-[cyclopropyl-(3-trifluoromethylbenzyl)carbamoyl]-3,9-diazabicyclo[3.3.1]non-6-en-3-yl)-3-hydroxy-5-oxopentanoic acid;

- 1:1-mixture of (*rac.*)-(3*R**)-5-((1*R**, 5*S**)-7-{4-[2-(2-chloro-4,5-dimethylphenoxy)ethoxy]phenyl}-6-{cyclopropyl-[2-(4-methoxyphenoxy)ethyl]-carbamoyl}-3,9-diazabicyclo[3.3.1]non-6-en-3-yl)-3-hydroxy-5-oxopentanoic acid and (*rac.*)-(3*S**)-5-((1*R**, 5*S**)-7-{4-[2-(2-chloro-4,5-dimethylphenoxy)ethoxy]phenyl}-6-{cyclopropyl-[2-(4-methoxyphenoxy)ethyl]carbamoyl}-3,9-diazabicyclo[3.3.1]non-6-en-3-yl)-3-hydroxy-5-oxopentanoic acid;
- 1:1-mixture of (*rac.*)-(3*R**)-5-[(1*R**, 5*S**)-7-{4-[2-(2-chloro-4,5-dimethylphenoxy)ethoxy]phenyl}-6-(cyclopropylphenethylcarbamoyl)-3,9-diazabicyclo[3.3.1]non-6-en-3-yl]-3-hydroxy-5-oxopentanoic acid and (*rac.*)-(3*S**)-5-[(1*R**, 5*S**)-7-{4-[2-(2-chloro-4,5-dimethylphenoxy)ethoxy]phenyl}-6-(cyclopropylphenethylcarbamoyl)-3,9-diazabicyclo[3.3.1]non-6-en-3-yl]-3-hydroxy-5-oxopentanoic acid;
- 1:1-mixture of (*rac.*)-(3*R**)-5-((1*R**, 5*S**)-7-{4-[2-(2-chloro-4,5-dimethylphenoxy)ethoxy]phenyl}-6-{[2-(2-chlorophenyl)ethyl]cyclopropylcarbamoyl}-3,9-diazabicyclo[3.3.1]non-6-en-3-yl)-3-hydroxy-5-oxopentanoic acid and (*rac.*)-(3*R**)-5-((1*R**, 5*S**)-7-{4-[2-(2-chloro-4,5-dimethylphenoxy)ethoxy]phenyl}-6-{[2-(2-chlorophenyl)ethyl]cyclopropylcarbamoyl}-3,9-diazabicyclo[3.3.1]non-6-en-3-yl)-3-hydroxy-5-oxopentanoic acid;
- 1:1-mixture of (*rac.*)-(3*R**)-5-((1*R**, 5*S**)-7-{4-[2-(2-chloro-4,5-dimethylphenoxy)ethoxy]phenyl}-6-{cyclopropyl-[2-(2,3-difluorophenyl)ethyl]-carbamoyl}-3,9-diazabicyclo[3.3.1]non-6-en-3-yl)-3-hydroxy-5-oxopentanoic acid and (*rac.*)-(3*S**)-5-((1*R**, 5*S**)-7-{4-[2-(2-chloro-4,5-dimethylphenoxy)ethoxy]phenyl}-6-{cyclopropyl-[2-(2,3-difluorophenyl)ethyl]carbamoyl}-3,9-diazabicyclo[3.3.1]non-6-en-3-yl)-3-hydroxy-5-oxopentanoic acid;
- 1:1-mixture of (*rac.*)-(3*R**)-5-((1*R**, 5*S**)-7-{4-[2-(2-chloro-4,5-dimethylphenoxy)ethoxy]phenyl}-6-{cyclopropyl-[2-(4-fluorophenyl)ethyl]carbamoyl}-3,9-diazabicyclo[3.3.1]non-6-en-3-yl)-3-hydroxy-5-oxopentanoic acid and (*rac.*)-(3*S**)-5-((1*R**, 5*S**)-7-{4-[2-(2-chloro-4,5-dimethylphenoxy)ethoxy]phenyl}-6-

{cyclopropyl-[2-(4-fluorophenyl)ethyl]carbamoyl}-3,9-diazabicyclo[3.3.1]non-6-en-3-yl)-3-hydroxy-5-oxopentanoic acid;

1:1-mixture of (*rac.*)-(3*R**)-5-{(*1R**, *5S**)-7-{4-[2-(2-chloro-4,5-dimethylphenoxy)ethoxy]phenyl}-6-[cyclopropyl(2-*o*-tolylethyl)carbamoyl]-3,9-diazabicyclo[3.3.1]non-6-en-3-yl}-3-hydroxy-5-oxopentanoic acid and (*rac.*)-(3*S**)-5-{(*1R**, *5S**)-7-{4-[2-(2-chloro-4,5-dimethylphenoxy)ethoxy]phenyl}-6-[cyclopropyl(2-*o*-tolylethyl)carbamoyl]-3,9-diazabicyclo[3.3.1]non-6-en-3-yl}-3-hydroxy-5-oxopentanoic acid;

10

1:1-mixture of (*rac.*)-(3*R**)-5-{(*1R**, *5S**)-7-{4-[2-(2-chloro-4,5-dimethylphenoxy)ethoxy]phenyl}-6-[cyclopropyl(2-*p*-tolylethyl)carbamoyl]-3,9-diazabicyclo[3.3.1]non-6-en-3-yl}-3-hydroxy-5-oxopentanoic acid and (*rac.*)-(3*S**)-5-{(*1R**, *5S**)-7-{4-[2-(2-chloro-4,5-dimethylphenoxy)ethoxy]phenyl}-6-[cyclopropyl(2-*p*-tolylethyl)carbamoyl]-3,9-diazabicyclo[3.3.1]non-6-en-3-yl}-3-hydroxy-5-oxopentanoic acid;

1:1-mixture of (*rac.*)-(3*R**)-5-{(*1R**, *5S**)-7-{4-[2-(2-chloro-4,5-dimethylphenoxy)ethoxy]phenyl}-6-[cyclopropyl-(3,5-dimethoxybenzyl)carbamoyl]-3,9-diazabicyclo[3.3.1]non-6-en-3-yl}-3-hydroxy-5-oxopentanoic acid and (*rac.*)-(3*S**)-5-{(*1R**, *5S**)-7-{4-[2-(2-chloro-4,5-dimethylphenoxy)ethoxy]phenyl}-6-[cyclopropyl-(3,5-dimethoxybenzyl)carbamoyl]-3,9-diazabicyclo[3.3.1]non-6-en-3-yl}-3-hydroxy-5-oxopentanoic acid;

25 1:1-mixture of (*1R*, *5S*)-7-{4-[2-(2-chloro-4,5-dimethylphenoxy)ethoxy]phenyl}-3-((2*S*, *4R*)-4-hydroxypyrrolidine-2-carbonyl)-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid (2-chlorobenzyl)cyclopropylamide and (*1S*, *5R*)-7-{4-[2-(2-chloro-4,5-dimethylphenoxy)ethoxy]phenyl}-3-((2*S*, *4R*)-4-hydroxypyrrolidine-2-carbonyl)-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid (2-chlorobenzyl)cyclopropylamide;

1:1-mixture of (*IR*, *5S*)-7-{4-[2-(2-chloro-4,5-dimethylphenoxy)ethoxy]phenyl}-3-((*2S*, *4R*)-4-hydroxypyrrolidine-2-carbonyl)-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid cyclopropyl-(2-methylbenzyl)amide and (*IS*, *5R*)-7-{4-[2-(2-chloro-4,5-dimethylphenoxy)ethoxy]phenyl}-3-((*2S*, *4R*)-4-hydroxypyrrolidine-2-carbonyl)-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid cyclopropyl-(2-methylbenzyl)amide;

(*rac.*)-(*IR**, *5S**)-3-acetyl-7-{4-[2-(2,3,5-trimethylphenoxy)ethyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid [2-(2-chlorophenyl)ethyl]-cyclopropylamide;

(*rac.*)-(*IR**, *5S**)-3-acetyl-7-{4-[2-(2,3,5-trimethylphenoxy)ethyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid cyclopropylphenethylamide;

(*rac.*)-(*IR**, *5S**)-3-acetyl-7-{4-[2-(2,3,5-trimethylphenoxy)ethyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid cyclopropyl-[2-(2,3-difluorophenyl)ethyl]amide;

(*rac.*)-(*IR**, *5S**)-3-acetyl-7-{4-[2-(2,3,5-trimethylphenoxy)ethyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid cyclopropyl-(3,5-dimethoxybenzyl)amide;

(*rac.*)-(*IR**, *5S**)-3-acetyl-7-{4-[2-(2,3,5-trimethylphenoxy)ethyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid cyclopropyl-(2-*p*-tolylethyl)amide;

(*rac.*)-(*IR**, *5S**)-3-acetyl-7-{4-[2-(2,3,5-trimethylphenoxy)ethyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid cyclopropyl-[2-(2-hydroxyethyl)benzyl]amide;

- (*rac.*)-(1*R*^{*}, 5*S*^{*})-3-acetyl-7-{4-[2-(2-chloro-4,5-dimethylphenoxy)ethoxy]-phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid cyclopropyl-[2-(2-hydroxyethyl)benzyl]amide;
- 5 (*rac.*)-5-((1*R*^{*}, 5*S*^{*})-6-(cyclopropylphenethylcarbamoyl)-7-{4-[2-(2,3,5-trimethylphenoxy)ethyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-en-3-yl)-5-oxopentanoic acid;
- 10 (*rac.*)-5-((1*R*^{*}, 5*S*^{*})-6-{[2-(2-chlorophenyl)ethyl]cyclopropylcarbamoyl}-7-{4-[2-(2,3,5-trimethylphenoxy)ethyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-en-3-yl)-5-oxopentanoic acid;
- 15 (*rac.*)-5-((1*R*^{*}, 5*S*^{*})-6-{cyclopropyl-[2-(2,3-difluorophenyl)ethyl]carbamoyl}-7-{4-[2-(2,3,5-trimethylphenoxy)ethyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-en-3-yl)-5-oxopentanoic acid;
- 20 (*rac.*)-5-((1*R*^{*}, 5*S*^{*})-6-{cyclopropyl-[2-(4-fluorophenyl)ethyl]carbamoyl}-7-{4-[2-(2,3,5-trimethyl-phenoxy)ethyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-en-3-yl)-5-oxopentanoic acid;
- 25 (*rac.*)-5-((1*R*^{*}, 5*S*^{*})-6-[cyclopropyl-(2-*o*-tolylethyl)carbamoyl]-7-{4-[2-(2,3,5-trimethylphenoxy)ethyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-en-3-yl)-5-oxo-pentanoic acid;
- 30 1:1-mixture of (*rac.*)-5-((1*R*^{*}, 5*S*^{*})-6-[cyclopropyl-((2*R*^{*})-2-hydroxy-2-phenylethyl)carbamoyl]-7-{4-[2-(2,3,5-trimethylphenoxy)ethyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-en-3-yl)-5-oxopentanoic acid and (*rac.*)-5-((1*R*^{*}, 5*S*^{*})-6-[cyclopropyl-((2*S*^{*})-2-hydroxy-2-phenylethyl)carbamoyl]-7-{4-[2-(2,3,5-trimethylphenoxy)ethyl]phenyl}-3,9-diazabicyclo-[3.3.1]non-6-en-3-yl)-5-oxopentanoic acid;

- (*rac.*)-5-((*1R*^{*}, *5S*^{*})-6-{cyclopropyl-[2-(2-hydroxyethyl)benzyl]carbamoyl}-7-{4-[2-(2,3,5-trimethylphenoxy)ethyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-en-3-yl)-5-oxo-pentanoic acid;
- 5 (*rac.*)-5-((*1R*^{*}, *5S*^{*})-7-{4-[3-(2-bromo-5-fluorophenoxy)propyl]phenyl}-6-{cyclopropyl-[2-(3-methoxyphenoxy)ethyl]carbamoyl}-3,9-diazabicyclo[3.3.1]-non-6-en-3-yl)-5-oxopentanoic acid;
- 10 (*rac.*)-5-((*1R*^{*}, *5S*^{*})-6-{[2-(2-chlorophenyl)ethyl]cyclopropylcarbamoyl}-7-{4-[2-(2,3,5-trimethylphenoxy)ethyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-en-3-yl)-2,2-dimethyl-5-oxopentanoic acid;
- 15 (*rac.*)-5-((*1R*^{*}, *5S*^{*})-6-{cyclopropyl-[2-(2,3-difluorophenyl)ethyl]carbamoyl}-7-{4-[2-(2,3,5-trimethylphenoxy)ethyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-en-3-yl)-2,2-dimethyl-5-oxopentanoic acid;
- 20 (*rac.*)-5-((*1R*^{*}, *5S*^{*})-6-{cyclopropyl-[2-(4-fluorophenyl)ethyl]carbamoyl}-7-{4-[2-(2,3,5-trimethylphenoxy)ethyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-en-3-yl)-2,2-dimethyl-5-oxopentanoic acid;
- 25 (*rac.*)-5-((*1R*^{*}, *5S*^{*})-6-[cyclopropyl-(2-*o*-tolylethyl)carbamoyl]-7-{4-[2-(2,3,5-trimethylphenoxy)ethyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-en-3-yl)-2,2-dimethyl-5-oxopentanoic acid;
- 30 (*rac.*)-5-((*1R*^{*}, *5S*^{*})-6-[cyclopropyl-(3,5-dimethoxybenzyl)carbamoyl]-7-{4-[2-(2,3,5-trimethylphenoxy)ethyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-en-3-yl)-2,2-dimethyl-5-oxopentanoic acid;

- (*rac.*)-5-((*1R*^{*}, *5S*^{*})-7-{4-[2-(2-chloro-4,5-dimethylphenoxy)ethoxy]phenyl}-6-{cyclopropyl-[2-(2-hydroxyethyl)benzyl]carbamoyl}-3,9-diazabicyclo[3.3.1]non-6-en-3-yl)-2,2-dimethyl-5-oxopentanoic acid;
- 5 (*rac.*)-(*IR*^{*}, *5S*^{*})-3-(4-carbamoylbutyryl)-7-{4-[2-(2,3,5-trimethylphenoxy)ethyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid cyclopropyl-phenethylamide;
- 10 (*rac.*)-(*IR*^{*}, *5S*^{*})-3-(4-carbamoylbutyryl)-7-{4-[2-(2,3,5-trimethylphenoxy)ethyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid [2-(2-chloro-phenyl)-ethyl]cyclopropylamide;
- 15 (*rac.*)-(*IR*^{*}, *5S*^{*})-3-(4-carbamoylbutyryl)-7-{4-[2-(2,3,5-trimethylphenoxy)ethyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid cyclopropyl-[2-(2,3-difluorophenyl)ethyl]amide;
- 20 (*rac.*)-(*IR*^{*}, *5S*^{*})-3-(4-carbamoylbutyryl)-7-{4-[2-(2,3,5-trimethylphenoxy)ethyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid cyclopropyl-[2-(4-fluorophenyl)ethyl]amide;
- 25 (*rac.*)-(*IR*^{*}, *5S*^{*})-3-(4-carbamoylbutyryl)-7-{4-[2-(2,3,5-trimethylphenoxy)ethyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid cyclopropyl-(3,5-dimethoxybenzyl)amide;
- 30 (*rac.*)-(*IR*^{*}, *5S*^{*})-3-(4-carbamoylbutyryl)-7-{4-[2-(2,3,5-trimethylphenoxy)ethyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid cyclopropyl-[2-(2-hydroxyethyl)benzyl]amide;

- (*rac.*)-(1*R*^{*}, 5*S*^{*})-3-(4-carbamoylbutyryl)-7-{4-[2-(2-chloro-4,5-dimethylphenoxy)ethoxy]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid cyclopropyl-[2-(2-hydroxyethyl)benzyl]amide;
- 5 (*rac.*)-acetic acid (1*R*^{*}, 5*S*^{*})-2-(2-[(3-acetyl-7-{4-[2-(2,3,5-trimethylphenoxy)ethyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carbonyl)cyclopropylamino]-methyl}phenyl)ethyl ester;
- 10 (*rac.*)-acetic acid (1*R*^{*}, 5*S*^{*})-2-(2-[(3-acetyl-7-{4-[2-(2-chloro-4,5-dimethylphenoxy)ethoxy]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carbonyl)-cyclopropylamino]methyl}phenyl)ethyl ester;
- 15 (*rac.*)-(1*R*^{*}, 5*S*^{*})-7-{4-[2-(2,3,5-trimethylphenoxy)ethyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid cyclopropyl-[2-(4-fluorophenyl)ethyl]-amide;
- 20 (*rac.*)-(1*R*^{*}, 5*S*^{*})-7-{4-[2-(2,3,5-trimethylphenoxy)ethyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid cyclopropyl-(3,5-dimethoxybenzyl)-amide;
- 25 1:1 mixture of (*rac.*)-5-[(1*R*^{*}, 5*S*^{*})-7-{4-[2-(2-chloro-4,5-dimethylphenoxy)ethoxy]phenyl}-6-[((2*R*^{*})2-hydroxy-2-phenylethyl)methylcarbamoyl]-3,9-diazabicyclo-[3.3.1]non-6-en-3-yl]-2,2-dimethyl-5-oxopentanoic acid and (*rac.*)-5-[(1*R*^{*}, 5*S*^{*})-7-{4-[2-(2-chloro-4,5-dimethylphenoxy)ethoxy]phenyl}-6-[((2*S*^{*})2-hydroxy-2-phenylethyl)methylcarbamoyl]-3,9-diazabicyclo-[3.3.1]non-6-en-3-yl]-2,2-dimethyl-5-oxopentanoic acid;
- 30

- (*rac.*)-(1*R*^{*}, 5*S*^{*})-3-acetyl-7-{4-[3-(2-bromo-5-fluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid (2-chlorobenzyl)cyclopropylamide;
- 5 (*rac.*)-(1*R*^{*}, 5*S*^{*})-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid benzylcyclopropylamide;
- 10 (*rac.*)-(1*R*^{*}, 5*S*^{*})-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid cyclopropyl-(2-fluorobenzyl)amide;
- 15 (*rac.*)-(1*R*^{*}, 5*S*^{*})-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid cyclopropyl-(3-trifluoromethylbenzyl)amide;
- 20 (*rac.*)-(1*R*^{*}, 5*S*^{*})-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid cyclopropyl-(2-methylbenzyl)amide;
- 25 (*rac.*)-(1*R*^{*}, 5*S*^{*})-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid cyclopropyl-[2-(4-methoxyphenoxy)-ethyl]amide;
- 30 (*rac.*)-(1*R*^{*}, 5*S*^{*})-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid [2-(2-chlorophenyl)ethyl]cyclopropylamide;

- (*rac.*)-(1*R*^{*}, 5*S*^{*})-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid cyclopropyl-(2-*p*-tolylethyl)amide;
- 5 (rac.)-(1*R*^{*}, 5*S*^{*})-3-acetyl-7-{4-[3-(2,5-difluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid (2-chlorobenzyl)cyclopropylamide;
- 10 (rac.)-(1*R*^{*}, 5*S*^{*})-3-acetyl-7-{4-[3-(2,3-dichlorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid (2-chlorobenzyl)cyclopropylamide;
- 15 (rac.)-(1*R*^{*}, 5*S*^{*})-3-acetyl-7-{4-[3-(2-chloro-5-fluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid (2-chlorobenzyl)cyclopropylamide;
- 20 (rac.)-(1*R*^{*}, 5*S*^{*})-3-acetyl-7-{4-[3-(3-cyanopyridin-2-yloxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid (2-chlorobenzyl)cyclopropylamide;
- 25 (rac.)-(1*R*^{*}, 5*S*^{*})-3-acetyl-7-{4-[2-(2,3-dimethylphenoxy)ethyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid (2-chlorobenzyl)cyclopropylamide;
- 30 (rac.)-(1*R*^{*}, 5*S*^{*})-3-acetyl-7-{4-[2-(2,5-dimethylphenoxy)ethyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid (2-chlorobenzyl)cyclopropylamide;
- (rac.)-(1*R*^{*}, 5*S*^{*})-3-acetyl-7-{4-[2-(2-chloro-4,5-dimethylphenoxy)ethyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid (2-chlorobenzyl)cyclopropylamide;

- (*rac.*)-(1*R*^{*,} 5*S*^{*})-3-acetyl-7-{4-[2-(2,4-dichlorophenoxy)ethyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid (2-chlorobenzyl)cyclopropylamide;
- 5 (rac.)-(1*R*^{*,} 5*S*^{*})-3-acetyl-7-{4-[2-(2,3-dichlorophenoxy)ethyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid (2-chlorobenzyl)cyclopropylamide;
- 10 (rac.)-(1*R*^{*,} 5*S*^{*})-3-acetyl-7-{4-[2-(2,6-difluorophenoxy)ethyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid (2-chlorobenzyl)cyclopropylamide;
- 15 (rac.)-(1*R*^{*,} 5*S*^{*})-3-acetyl-7-{4-[2-(3,5-dichlorophenoxy)ethyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid (2-chlorobenzyl)cyclopropylamide;
- 20 (rac.)-(1*R*^{*,} 5*S*^{*})-3-acetyl-7-{4-[2-(2-chloro-5-methylphenoxy)ethyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid (2-chlorobenzyl)cyclopropylamide;
- 25 (rac.)-(1*R*^{*,} 5*S*^{*})-3-acetyl-7-{4-[2-(2-chloro-5-fluorophenoxy)ethyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid (2-chlorobenzyl)cyclopropylamide;
- 30 (rac.)-(1*R*^{*,} 5*S*^{*})-3-acetyl-7-{4-[2-(2,3,6-trichlorophenoxy)ethyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid (2-chlorobenzyl)cyclopropylamide;
- (rac.)-(1*R*^{*,} 5*S*^{*})-3-acetyl-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid cyclopropyl-(2-methoxybenzyl)-amide;

- (*rac.*)-(1*R*^{*}, 5*S*^{*})-3-acetyl-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid cyclopropyl-(5-fluoro-2-methoxybenzyl)amide;
- 5 (*rac.*)-(1*R*^{*}, 5*S*^{*})-3-acetyl-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid cyclopropyl-(2,6-difluorobenzyl)amide;
- 10 (*rac.*)-(1*R*^{*}, 5*S*^{*})-3-acetyl-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid cyclopropyl-(3-fluoro-2-methylbenzyl)amide;
- 15 (*rac.*)-(1*R*^{*}, 5*S*^{*})-3-acetyl-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid (2-chloro-3,6-difluorobenzyl)cyclopropylamide;
- 20 (*rac.*)-(1*R*^{*}, 5*S*^{*})-3-acetyl-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid cyclopropyl-(2,3-difluorobenzyl)amide;
- 25 (*rac.*)-(1*R*^{*}, 5*S*^{*})-3-acetyl-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid cyclopropyl-(4-fluorobenzyl)amide;
- 30 (*rac.*)-(1*R*^{*}, 5*S*^{*})-7-{4-[3-(2-bromo-5-fluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid (2-chlorobenzyl)ethylamide;
- (*rac.*)-(1*R*^{*}, 5*S*^{*})-7-{4-[3-(2-bromo-5-fluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid cyclopropyl-(3,5-dimethoxybenzyl)amide;

- (*rac.*)-(1*R*^{*, 5S*})-7-{4-[3-(2-bromo-5-fluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid cyclopropyl-(2-methylbenzyl)amide;
- (*rac.*)-(1*R*)-7-{4-[3-(2-bromo-5-fluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid cyclopropyl-[2-(4-methoxyphenoxy)-ethyl]amide;
- (*rac.*)-(1*R*)-3-methyl-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid (2-chlorobenzyl)cyclopropylamide;
- (*rac.*)-(1*R*)-3-ethyl-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid (2-chlorobenzyl)cyclopropylamide;
- (*rac.*)-(1*R*)-3-(2-aminoacetyl)-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid (2-chlorobenzyl)cyclopropylamide;
- (*rac.*)-(1*R*)-3-(3-aminopropionyl)-7-{4-[3-(2,3,6-trifluorophenoxy)-propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid (2-chlorobenzyl)cyclopropylamide;
- (*rac.*)-(1*R*)-3-acetyl-7-{4-[2-(2,4,5-trichlorophenoxy)ethoxy]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid (2-chlorobenzyl)cyclopropylamide;
- (*rac.*)-(1*R*)-3-acetyl-7-{4-[2-(2-bromo-5-fluorophenoxy)ethoxy]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid (2-chloro-3-trifluoromethylbenzyl)cyclopropylamide;

- (*rac.*)-(1*R*^{*}, 5*S*^{*})-3-acetyl-7-{4-[2-(2,4,5-trichlorophenoxy)ethoxy]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid (2-bromobenzyl)cyclopropylamide;
- 5 (*rac.*)-(1*R*^{*}, 5*S*^{*})-3-acetyl-7-{4-[2-(2-bromo-5-fluorophenoxy)ethoxy]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid cyclopropyl-(2,3-dimethylbenzyl)amide;
- 10 (*rac.*)-(1*R*^{*}, 5*S*^{*})-3-acetyl-7-{4-[2-(2,3-dichlorophenoxy)ethoxy]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid (2-chloro-3-trifluoromethylbenzyl)cyclopropylamide;
- 15 (*rac.*)-(1*R*^{*}, 5*S*^{*})-3-acetyl-7-{4-[2-(2-chloro-4,5-dimethylphenoxy)ethoxy]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid (3-chlorobenzyl)cyclopropylamide;
- 20 (*rac.*)-(1*R*^{*}, 5*S*^{*})-3-acetyl-7-{4-[2-(2,4,5-trichlorophenoxy)ethoxy]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid (2-chlorobenzyl)ethylamide;
- 25 (*rac.*)-(1*R*^{*}, 5*S*^{*})-3-acetyl-7-{4-[2-(2-chloro-5-fluorophenoxy)ethoxy]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid (2-bromobenzyl)cyclopropylamide;
- 30 (*rac.*)-(1*R*^{*}, 5*S*^{*})-3-acetyl-7-{4-[2-(4-chloro-2-methylphenoxy)ethoxy]-phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid cyclopropyl-(2,3-dichlorobenzyl)amide;
- 35 (*rac.*)-(1*R*^{*}, 5*S*^{*})-3-acetyl-7-{4-[2-(2-chloro-4,5-dimethylphenoxy)ethoxy]-phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid cyclopropyl-(3-methoxybenzyl)amide;

- (*rac.*)-(1*R*^{*}, 5*S*^{*})-3-acetyl-7-{4-[2-(2-bromo-5-fluorophenoxy)ethoxy]-phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid (2-bromobenzyl)-cyclopropylamide;
- 5 (*rac.*)-(1*R*^{*}, 5*S*^{*})-3-acetyl-7-{4-[2-(4-chloro-2-methylphenoxy)ethoxy]-phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid cyclopropyl-(2,3-dimethylbenzyl)amide;
- 10 (*rac.*)-(1*R*^{*}, 5*S*^{*})-3-acetyl-7-{4-[2-(2,4,5-trichlorophenoxy)ethoxy]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid cyclopropyl-(3-methoxybenzyl)amide;
- 15 (*rac.*)-(1*R*^{*}, 5*S*^{*})-3-acetyl-7-{4-[2-(2-chloro-4,5-dimethylphenoxy)ethoxy]-phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid cyclopropyl-(2-fluoro-5-methoxybenzyl)amide;
- 20 (*rac.*)-(1*R*^{*}, 5*S*^{*})-3-acetyl-7-{4-[2-(4-chloro-2-methylphenoxy)ethoxy]-phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid (2-bromobenzyl)-cyclopropylamide;
- 25 (*rac.*)-(1*R*^{*}, 5*S*^{*})-3-acetyl-7-{4-[2-(2-chloro-5-fluorophenoxy)ethoxy]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid (2-chlorobenzyl)-cyclopropylamide;
- 30 (*rac.*)-(1*R*^{*}, 5*S*^{*})-3-Acetyl-7-{4-[2-(2,4,5-trichlorophenoxy)ethoxy]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid cyclopropyl-(3,5-dimethoxybenzyl)amide;
- 35 (*rac.*)-(1*R*^{*}, 5*S*^{*})-3-acetyl-7-{4-[2-(2,3-dichlorophenoxy)ethoxy]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid (2-bromobenzyl)cyclopropylamide;

- (*rac.*)-(1*R*^{*}, 5*S*^{*})-3-acetyl-7-{4-[2-(2-chloro-4,5-dimethylphenoxy)ethoxy]-phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid (6-chloro-benzo[1,3]dioxol-5-ylmethyl)-cyclopropylamide;
- 5 (*rac.*)-(1*R*^{*}, 5*S*^{*})-3-acetyl-7-{4-[2-(2,3-dichlorophenoxy)ethoxy]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid (2-chlorobenzyl)cyclopropylamide;
- 10 (*rac.*)-(1*R*^{*}, 5*S*^{*})-3-acetyl-7-{4-[2-(4-chloro-2-methylphenoxy)ethoxy]-phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid (2-chloro-3-trifluoromethylbenzyl)cyclopropylamide;
- 15 (*rac.*)-(1*R*^{*}, 5*S*^{*})-3-acetyl-7-{4-[2-(2-chloro-4,5-dimethylphenoxy)ethoxy]-phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid cyclopropyl-(3-methylbenzyl)amide;
- 20 (*rac.*)-(1*R*^{*}, 5*S*^{*})-3-acetyl-7-{4-[2-(2,3-dichlorophenoxy)ethoxy]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid (2-chlorobenzyl)ethylamide;
- 25 (*rac.*)-(1*R*^{*}, 5*S*^{*})-3-acetyl-7-{4-[2-(2-chloro-5-fluorophenoxy)ethoxy]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid cyclopropyl-(2-fluoro-5-methoxybenzyl)amide;
- 30 (*rac.*)-(1*R*^{*}, 5*S*^{*})-3-acetyl-7-{4-[2-(4-chloro-2-methylphenoxy)ethoxy]-phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid (2-chlorobenzyl)-cyclopropylamide;
- (*rac.*)-(1*R*^{*}, 5*S*^{*})-3-acetyl-7-{4-[2-(2-bromo-5-fluorophenoxy)ethoxy]-phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid cyclopropyl-(3,5-dimethoxybenzyl)amide;

- (*rac.*)-(1*R*^{*}, 5*S*^{*})-3-acetyl-7-{4-[2-(2-chloro-5-fluorophenoxy)ethoxy]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid cyclopropyl-(3-methoxybenzyl)amide;
- 5 (*rac.*)-(1*R*^{*}, 5*S*^{*})-3-acetyl-7-{4-[2-(2,4,5-trichlorophenoxy)ethoxy]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid (3-chlorobenzyl)cyclopropylamide;
- 10 (*rac.*)-(1*R*^{*}, 5*S*^{*})-3-acetyl-7-{4-[2-(2,6-dichloro-4-methylphenoxy-ethoxy]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid cyclopropyl-(3,5-difluorobenzyl)amide;
- 15 (*rac.*)-(1*R*^{*}, 5*S*^{*})-3-acetyl-7-{4-[2-(2,4,5-trichlorophenoxy)ethoxy]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid cyclopropyl-(2-fluoro-5-methoxybenzyl)amide;
- 20 (*rac.*)-(1*R*^{*}, 5*S*^{*})-3-acetyl-7-{4-[2-(2,4,5-trichlorophenoxy)ethoxy]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid cyclopropyl-(3,4-dimethoxybenzyl)amide;
- 25 (*rac.*)-(1*R*^{*}, 5*S*^{*})-3-acetyl-7-{4-[2-(2-chloro-5-fluorophenoxy)ethoxy]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid (6-chloro-benzo[1,3]-dioxol-5-ylmethyl)cyclopropylamide;
- 30 (*rac.*)-(1*R*^{*}, 5*S*^{*})-3-Acetyl-7-{4-[2-(2-bromo-5-fluorophenoxy)ethoxy]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid (3-chlorobenzyl)cyclopropylamide;
- 35 (*rac.*)-(1*R*^{*}, 5*S*^{*})-3-Acetyl-7-{4-[2-(2,3-dichlorophenoxy)ethoxy]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid cyclopropyl-(3,4-dimethoxybenzyl)amide;

- (*rac.*)-(1*R*^{*}, 5*S*^{*})-3-acetyl-7-{4-[2-(2-chloro-5-fluorophenoxy)ethoxy]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid (2-chlorobenzyl)-ethylamide;
- 5 (rac.)-(1*R*^{*}, 5*S*^{*})-3-acetyl-7-{4-[2-(4-chloro-2-methylphenoxy)ethoxy]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid cyclopropyl-(3,5-dimethoxybenzyl)amide;
- 10 (rac.)-(1*R*^{*}, 5*S*^{*})-3-acetyl-7-{4-[2-(2,4,5-trichlorophenoxy)ethoxy]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid cyclopropyl-(3-methylbenzyl)amide;
- 15 (rac.)-(1*R*^{*}, 5*S*^{*})-3-acetyl-7-{4-[2-(2,3-dichlorophenoxy)ethoxy]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid cyclopropyl-(2-fluoro-5-methoxybenzyl)amide;
- 20 (rac.)-(1*R*^{*}, 5*S*^{*})-3-acetyl-7-{4-[2-(2-chloro-4,5-dimethylphenoxy)ethoxy]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid cyclopropyl-(3,5-difluorobenzyl)amide;
- 25 (rac.)-(1*R*^{*}, 5*S*^{*})-3-acetyl-7-{4-[2-(2,3-dichlorophenoxy)ethoxy]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid cyclopropyl-(3,5-dimethoxybenzyl)amide;
- 30 (rac.)-(1*R*^{*}, 5*S*^{*})-3-acetyl-7-{4-[2-(2-chloro-5-fluorophenoxy)ethoxy]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid (2-chlorobenzyl)-cyclopropylamide;
- (rac.)-(1*R*^{*}, 5*S*^{*})-3-acetyl-7-{4-[2-(2-bromo-5-fluorophenoxy)ethoxy]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid (2-chlorobenzyl)-cyclopropylamide;

- (*rac.*)-(1*R*^{*}, 5*S*^{*})-3-acetyl-7-{4-[2-(2-chloro-4,5-dimethylphenoxy)ethoxy]-phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid cyclopropyl-(3,4-dimethoxybenzyl)amide;
- 5 (*rac.*)-(1*R*^{*}, 5*S*^{*})-3-acetyl-7-{4-[2-(2-chloro-4,5-dimethylphenoxy)ethoxy]-phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid cyclopropyl-(3-trifluoromethoxybenzyl)amide;
- 10 (*rac.*)-(1*R*^{*}, 5*S*^{*})-3-acetyl-7-{4-[2-(2-chloro-5-fluorophenoxy)ethoxy]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid cyclopropyl-(3-methylbenzyl)amide;
- 15 (*rac.*)-(1*R*^{*}, 5*S*^{*})-3-acetyl-7-{4-[2-(2-chloro-5-fluorophenoxy)ethoxy]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid cyclopropyl-(3,5-dimethoxybenzyl)amide;
- 20 (*rac.*)-(1*R*^{*}, 5*S*^{*})-3-acetyl-7-{4-[2-(2,3-dichlorophenoxy)ethoxy]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid cyclopropyl-(3-methylbenzyl)amide;
- 25 (*rac.*)-(1*R*^{*}, 5*S*^{*})-3-acetyl-7-{4-[2-(2-bromo-5-fluorophenoxy)ethoxy]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid cyclopropyl-(2-fluoro-5-methoxybenzyl)amide;
- 30 (*rac.*)-(1*R*^{*}, 5*S*^{*})-3-acetyl-7-{4-[2-(2-bromo-5-fluorophenoxy)ethoxy]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid (2-chlorobenzyl)ethylamide;
- (*rac.*)-(1*R*^{*}, 5*S*^{*})-3-acetyl-7-{4-[2-(2,3-dichlorophenoxy)ethoxy]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid (6-chlorobenzo[1,3]dioxol-5-ylmethyl)cyclopropylamide;

- (*rac.*)-(1*R*^{*,} 5*S*^{*})-3-acetyl-7-{4-[2-(2,3-dichlorophenoxy)ethoxy]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid (3-chlorobenzyl)cyclopropylamide;
- 5 (*rac.*)-(1*R*^{*,} 5*S*^{*})-3-acetyl-7-{4-[2-(4-chloro-2-methylphenoxy)ethoxy]-phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid (6-chloro-benzo[1,3]dioxol-5-ylmethyl)cyclopropylamide;
- 10 (*rac.*)-(1*R*^{*,} 5*S*^{*})-3-acetyl-7-{4-[2-(2-bromo-5-fluorophenoxy)ethoxy]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid cyclopropyl-(3-methoxybenzyl)amide;
- 15 (*rac.*)-(1*R*^{*,} 5*S*^{*})-3-acetyl-7-{4-[2-(4-chloro-2-methylphenoxy)ethoxy]-phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid (2-chlorobenzyl)ethylamide;
- 20 (*rac.*)-(1*R*^{*,} 5*S*^{*})-3-acetyl-7-{4-[2-(4-chloro-2-methylphenoxy)ethoxy]-phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid cyclopropyl-(3-methylbenzyl)amide;
- 25 (*rac.*)-(1*R*^{*,} 5*S*^{*})-3-acetyl-7-{4-[2,(2,4,5-trichlorophenoxy)ethoxy]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid (6-chlorobenzo[1,3]dioxol-5-ylmethyl)cyclopropylamide;
- 30 (*rac.*)-(1*R*^{*,} 5*S*^{*})-3-acetyl-7-{4-[3-(2,6-dichloro-4-methylphenoxy)propyl]-phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid (2-chlorobenzyl)-cyclopropylamide formate salt;
- 35 (*rac.*)-(1*R*^{*,} 5*S*^{*})-7-{4-[3-(2-bromo-5-fluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid (2-chlorobenzyl)cyclopropylamide.

Most preferred compounds of general formula I are those selected from the group consisting of

- (*rac.*)-(2-methoxyphenyl)acetic acid (*1R*^{*,} *5S*^{*})-7-{4-[3-(2-methoxybenzyloxy)-
5 propoxy]phenyl}-3-(2-thiophen-2-ylacetyl)-3,9-diazabicyclo[3.3.1]non-6-en-6-yl-
methyl ester;
- (*rac.*)-(2-methoxyphenyl)acetic acid (*1R*^{*,} *5S*^{*})-3-[2-(4-chlorophenyl)acetyl]-7-
10 {4-[3-(2-methoxybenzyloxy)propoxy]phenyl}-3,9-diazabicyclo[3.3.1]non-6-en-6-
ylmethyl ester;
- (*rac.*)-(2-methoxyphenyl)acetic acid (*1R*^{*,} *5S*^{*})-7-{4-[3-(2-methoxybenzyloxy)-
propoxy]phenyl}-3-(quinoxaline-2-carbonyl)-3,9-diazabicyclo[3.3.1]non-6-en-6-
ylmethyl ester;
- 15 (*rac.*)-(*IR*^{*,} *5S*^{*})-3-acetyl-7-{4-[3-(2-methoxybenzyloxy)propoxy]phenyl}-3,9-di-
azabicyclo[3.3.1]non-6-ene-6-carboxylic acid [2-(2-chlorophenyl)ethyl]methyl-
amide;
- 20 (*rac.*)-(2-methoxyphenyl)acetic acid (*1R*^{*,} *5S*^{*})-3-(benzo[b]thiophene-3-
carbonyl)-7-{4-[3-(2-methoxybenzyloxy)propoxy]phenyl}-3,9-diazabicyclo-
[3.3.1]non-6-en-6-yl-methyl ester;
- (*rac.*)-(2-methoxyphenyl)acetic acid (*1R*^{*,} *5S*^{*})-3-acetyl-7-{4-[3-(2-methoxy-
25 benzyloxy)propoxy]phenyl}-3,9-diazabicyclo[3.3.1]non-6-en-6-ylmethyl ester;
- (*rac.*)-(2-methoxyphenyl)acetic acid (*1R*^{*,} *5S*^{*})-7-{4-[3-(2-methoxybenzyloxy)-
propoxy]phenyl}-3-phenylmethanesulfonyl-3,9-diazabicyclo[3.3.1]non-6-en-6-
ylmethyl ester;

- (*rac.*)-(1*R*^{*,} 5*S*^{*})-3-acetyl-7-{4-[3-(2-methoxybenzyloxy)propoxy]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid [2-(4-methoxyphenyl)ethyl]-methylamide;
- 5 (*rac.*)-(1*R*^{*,} 5*S*^{*})-3-acetyl-7-{4-[2-(2-chloro-5-methylphenoxy)ethyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid methylphenethylamide;
- (*rac.*)-(1*R*^{*,} 5*S*^{*})-3-acetyl-7-{4-[2-(2-bromo-5-fluorophenoxy)ethyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid methylphenethylamide;
- 10 (*rac.*)-(2-methoxyphenyl)acetic acid (1*R*^{*,} 5*S*^{*})-3-[2-(4-chlorophenyl)acetyl]-7-{6-[3-(2-methoxybenzyloxy)propoxy]pyridin-3-yl}-3,9-diazabicyclo[3.3.1]non-6-en-6-ylmethyl ester;
- 15 (*rac.*)-(1*R*^{*,} 5*S*^{*})-3-acetyl-7-{4-[2-(2,5-dimethylphenoxy)ethyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid methylphenethylamide;
- (*rac.*)-(1*R*^{*,} 5*S*^{*})-3-acetyl-7-{4-[3-(2-methoxybenzyloxy)propoxy]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid [2-(4-chlorophenyl)ethyl]methylamide;
- 20 (*rac.*)-(1*R*^{*,} 5*S*^{*})-3-acetyl-7-{4-[3-(2-methoxybenzyloxy)propoxy]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid [2-(3-chlorophenyl)ethyl]methylamide;
- 25 (*rac.*)-(1*R*^{*,} 5*S*^{*})-3-acetyl-7-{4-[3-(2-methoxybenzyloxy)propoxy]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid ethylphenethylamide;
- (*rac.*)-(1*R*^{*,} 5*S*^{*})-3-acetyl-7-{4-[3-(2-methoxybenzyloxy)propoxy]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid [2-(3-methoxyphenyl)ethyl]methylamide;
- 30 (*rac.*)-(1*R*^{*,} 5*S*^{*})-3-acetyl-7-{4-[3-(2-methoxybenzyloxy)propoxy]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid [2-(3-methoxyphenyl)ethyl]methylamide;

- (*rac.*)-(1*R*^{*,} 5*S*^{*})-3-acetyl-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-di-azabicyclo[3.3.1]non-6-ene-6-carboxylic acid methylphenethylamide;
- 5 (rac.)-(1*R*^{*,} 5*S*^{*})-3-acetyl-7-{4-[3-(2-methoxybenzyloxy)propoxy]phenyl}-3,9-di-azabicyclo[3.3.1]non-6-ene-6-carboxylic acid methylphenethylamide;
- 10 (rac.)-(1*R*^{*,} 5*S*^{*})-3-acetyl-7-{4-[3-(2-bromo-5-fluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid methylphenethylamide;
- 15 (rac.)-(2-methoxyphenyl)acetic acid (1*R*^{*,} 5*S*^{*})-7-{4-[3-(2-methoxybenzyloxy)propoxy]phenyl}-3-methyl-3,9-diazabicyclo[3.3.1]non-6-en-6-ylmethyl ester;
- 20 (rac.)-(1*R*^{*,} 5*S*^{*})-3-acetyl-7-{4-[3-(2-methoxybenzyloxy)propoxy]phenyl}-3,9-di-azabicyclo[3.3.1]non-6-ene-6-carboxylic acid [2-(3,4-dimethoxyphenyl)ethyl]-methylamide;
- 25 (rac.)-(1*R*^{*,} 5*S*^{*})-3-acetyl-7-{4-[2-(2,3,5-trimethylphenoxy)ethyl]phenyl}-3,9-di-azabicyclo[3.3.1]non-6-ene-6-carboxylic acid methylphenethylamide;
- 30 (rac.)-(1*R*^{*,} 5*S*^{*})-3-acetyl-7-{4-[3-(2,5-dichlorophenoxy)propyl]phenyl}-3,9-di-azabicyclo[3.3.1]non-6-ene-6-carboxylic acid methylphenethylamide;
- (rac.)-N-((1*R*^{*,} 5*S*^{*})-3-acetyl-7-{4-[2-(2-chloro-4,5-dimethylphenoxy)ethyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid methylphenethylamide;
- (rac.)-(2-methoxyphenyl)acetic acid (1*R*^{*,} 5*S*^{*})-7-{4-[3-(2-methoxybenzyloxy)propoxy]phenyl}-3,9-diazabicyclo[3.3.1]non-6-en-6-ylmethyl ester;

- (*rac.*)-*N*-((*1R*^{*}, *5S*^{*})-3-acetyl-7-{4-[3-(2-methoxybenzyloxy)propoxy]phenyl}-3,9-diazabicyclo[3.3.1]non-6-en-6-ylmethyl)-2-(2-methoxyphenyl)-*N*-methylacetamide;
- 5 (*rac.*)-*N*-((*1R*^{*}, *5S*^{*})-3-acetyl-7-{4-[2-(2-*tert*-butyl-4-methylphenoxy)ethyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid methylphenethylamide;
- (*rac.*)-(*1R*^{*}, *5S*^{*})-3-acetyl-7-{4-[2-(2,5-difluorophenoxy)ethyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid methylphenethylamide;
- 10 (*rac.*)-(*1R*^{*}, *5S*^{*})-3-acetyl-7-{4-[3-(2-methoxybenzyloxy)propoxy]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid methyl(3-phenylpropyl)amide;
- 15 (*rac.*)-(*1R*^{*}, *5S*^{*})-3-acetyl-7-{4-[3-(2,3-dichlorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid methylphenethylamide;
- (*rac.*)-(*1R*^{*}, *5S*^{*})-3-acetyl-7-{4-[3-(2-acetylphenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid methylphenethylamide;
- 20 (*rac.*)-(*1R*^{*}, *5S*^{*})-3-acetyl-7-{4-[3-(2-methoxybenzyloxy)propoxy]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid [2-(2-methoxyphenyl)ethyl]methylamide;
- 25 (*rac.*)-(*1R*^{*}, *5S*^{*})-3-acetyl-7-{4-[3-(2-methoxybenzyloxy)propoxy]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid benzylmethyleamide;
- (*rac.*)-(*1R*^{*}, *5S*^{*})-3-acetyl-7-{4-[3-(2,5-difluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid methylphenethylamide;
- 30 (*rac.*)-(*1R*^{*}, *5S*^{*})-3-acetyl-7-[4-(2-*o*-tolyloxyethyl)phenyl]-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid methylphenethylamide;

- (*rac.*)-(1*R*^{*,} 5*S*^{*})-3-acetyl-7-{4-[2-(3-isopropylphenoxy)ethyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid methylphenethylamide;
- (*rac.*)-(1*R*^{*,} 5*S*^{*})-3-acetyl-7-{4-[3-(2-chlorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid (2-chlorobenzyl)cyclopropylamide;
- (*rac.*)-(1*R*^{*,} 5*S*^{*})-5-[7-{4-[2-(2-bromo-5-fluorophenoxy)ethyl]phenyl}-6-(methylphenethylcarbamoyl)-3,9-diazabicyclo[3.3.1]non-6-en-3-yl]-5-oxopentanoic acid;
- 10 (*rac.*)-(1*R*^{*,} 5*S*^{*})-3-acetyl-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid (2-chlorobenzyl)cyclopropylamide;
- 15 (*rac.*)-(1*R*^{*,} 5*S*^{*})-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid (2-chlorobenzyl)cyclopropylamide;
- (*rac.*)-5-((1*R*^{*,} 5*S*^{*})-6-[(2-chlorobenzyl)cyclopropylcarbamoyl]-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-en-3-yl)-5-oxopentanoic acid;
- 20 (*rac.*)-(1*R*^{*,} 5*S*^{*})-6-[(2-chlorobenzyl)cyclopropylcarbamoyl]-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-9-carboxylic acid 2,2,2-trichloro-1,1-dimethylethyl ester;
- 25 (*rac.*)-5-((1*R*^{*,} 5*S*^{*})-6-[(2-chlorobenzyl)cyclopropylcarbamoyl]-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-en-3-yl)-2,2-dimethyl-5-oxopentanoic acid;
- 30 (*rac.*)-5-((1*R*^{*,} 5*S*^{*})-6-[(2-chlorobenzyl)cyclopropylcarbamoyl]-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-en-3-yl)-5-oxopentanoic acid methyl ester;

- 1:1-mixture of (*1R, 5S*)-3-((*1S, 4R*)-4-hydroxypyrrolidine-2-carbonyl)-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid (2-chlorobenzyl)cyclopropylamide and (*1S, 5R*)-3-((*1S, 4R*)-4-hydroxypyrrolidine-2-carbonyl)-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-5 3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid (2-chlorobenzyl)cyclopropylamide;
- (*rac.*)-(*1R**, *5S**)-3-(4-carbamoylbutyryl)-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid (2-chlorobenzyl)cyclopropylamide; 10
- (*rac.*)-(*1R**, *5S**)-3-(4-carbamoylbutyryl)-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid benzylcyclopropylamide; 15
- (*rac.*)-(*1R**, *5S**)-3-(4-carbamoylbutyryl)-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid (2-chlorobenzyl)ethylamide;
- 20 (*rac.*)-(*1R**, *5S**)-3-(4-carbamoylbutyryl)-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid cyclopropyl-(2-fluorobenzyl)amide;
- (*rac.*)-(*1R**, *5S**)-3-(4-carbamoylbutyryl)-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid cyclopropyl-(3-trifluoromethylbenzyl)amide; 25
- (*rac.*)-(*1R**, *5S**)-3-(4-carbamoylbutyryl)-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid cyclopropyl-(2-methylbenzyl)amide; 30

- (*rac.*)-(1*R*^{*,} 5*S*^{*})-3-(4-carbamoylbutyryl)-7-{4-[3-(2,3,6-trifluorophenoxy)-propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid cyclopropyl-phenethylamide;
- 5 (*rac.*)-(1*R*^{*,} 5*S*^{*})-3-(4-carbamoylbutyryl)-7-{4-[3-(2,3,6-trifluorophenoxy)-propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid [2-(2-chlorophenyl)ethyl]cyclopropylamide;
- 10 (*rac.*)-(1*R*^{*,} 5*S*^{*})-3-(4-carbamoylbutyryl)-7-{4-[3-(2,3,6-trifluorophenoxy)-propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid cyclopropyl-(3,5-dimethoxybenzyl)amide;
- 15 1:1-mixture of (*rac.*)-(3*R*^{*})-5-((1*R*^{*,} 5*S*^{*})-6-[(2-chlorobenzyl)cyclopropylcarbamoyl]-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-en-3-yl)-3-hydroxy-5-oxopentanoic acid and (*rac.*)-(3*S*^{*})-5-((1*R*^{*,} 5*S*^{*})-6-[(2-chlorobenzyl)cyclopropylcarbamoyl]-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-en-3-yl)-3-hydroxy-5-oxopentanoic acid;
- 20 20 1:1-mixture of (*rac.*)-(3*R*^{*})-5-((1*R*^{*,} 5*S*^{*})-6-(benzylcyclopropylcarbamoyl)-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-en-3-yl)-3-hydroxy-5-oxopentanoic acid and (*rac.*)-(3*S*^{*})-5-((1*R*^{*,} 5*S*^{*})-6-(benzylcyclopropylcarbamoyl)-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-en-3-yl)-3-hydroxy-5-oxopentanoic acid;
- 25 25 1:1-mixture of (*rac.*)-(3*R*^{*})-5-((1*R*^{*,} 5*S*^{*})-6-[(2-chlorobenzyl)ethylcarbamoyl]-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-en-3-yl)-3-hydroxy-5-oxopentanoic acid and (*rac.*)-(3*S*^{*})-5-((1*R*^{*,} 5*S*^{*})-6-[(2-chlorobenzyl)ethylcarbamoyl]-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-en-3-yl)-3-hydroxy-5-oxopentanoic acid;
- 30 30 3,9-diazabicyclo[3.3.1]non-6-en-3-yl)-3-hydroxy-5-oxopentanoic acid;

- 1:1-mixture of (*rac.*)-(3*R**)-5-((1*R**, 5*S**)-6-[cyclopropyl-(2-fluorobenzyl)-carbamoyl]-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo-[3.3.1]non-6-en-3-yl)-3-hydroxy-5-oxopentanoic acid and (*rac.*)-(3*S**)-5-((1*R**, 5*S**)-6-[cyclopropyl-(2-fluorobenzyl)carbamoyl]-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo-[3.3.1]non-6-en-3-yl)-3-hydroxy-5-oxopentanoic acid;
- 1:1-mixture of (*rac.*)-(3*R**)-5-((1*R**, 5*S**)-6-[cyclopropyl-(3-trifluoromethylbenzyl)carbamoyl]-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-en-3-yl)-3-hydroxy-5-oxopentanoic acid and (*rac.*)-(3*S**)-5-((1*R**, 5*S**)-6-[cyclopropyl-(3-trifluoromethylbenzyl)carbamoyl]-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]-phenyl}-3,9-diazabicyclo[3.3.1]non-6-en-3-yl)-3-hydroxy-5-oxopentanoic acid;
- 1:1-mixture of (*rac.*)-(3*R**)-5-((1*R**, 5*S**)-6-[cyclopropyl-(2-methylbenzyl)-carbamoyl]-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo-[3.3.1]non-6-en-3-yl)-3-hydroxy-5-oxopentanoic acid and (*rac.*)-(3*S**)-5-((1*R**, 5*S**)-6-[cyclopropyl-(2-methylbenzyl)carbamoyl]-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-en-3-yl)-3-hydroxy-5-oxopentanoic acid;
- 1:1-mixture of (*rac.*)-(3*R**)-5-((1*R**, 5*S**)-6-[cyclopropyl-[2-(4-methoxyphenoxy)ethyl]carbamoyl]-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-en-3-yl)-3-hydroxy-5-oxopentanoic acid and (*rac.*)-(3*S**)-5-((1*R**, 5*S**)-6-[cyclopropyl-[2-(4-methoxyphenoxy)ethyl]carbamoyl]-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-en-3-yl)-3-hydroxy-5-oxopentanoic acid;
- 1:1-mixture of (*rac.*)-(3*R**)-5-((1*R**, 5*S**)-6-[cyclopropyl-(2-*m*-tolyloxyethyl)-carbamoyl]-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo-[3.3.1]non-6-en-3-yl)-3-hydroxy-5-oxopentanoic acid and (*rac.*)-(3*S**)-5-((1*R**, 5*S**)-6-[cyclopropyl-(2-*m*-tolyloxyethyl)carbamoyl]-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-en-3-yl)-3-hydroxy-5-oxopentanoic acid;

phenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-en-3-yl)-3-hydroxy-5-oxopentanoic acid;

1:1-mixture of (*rac.*)-(3*R**)-5-((*1R**, 5*S**)-6-{[2-(2-chlorophenyl)ethyl]-cyclopropylcarbamoyl}-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-en-3-yl)-3-hydroxy-5-oxopentanoic acid and (*rac.*)-(3*S**)-5-((*1R**, 5*S**)-6-{[2-(2-chlorophenyl)ethyl]cyclopropylcarbamoyl}-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-en-3-yl)-3-hydroxy-5-oxopentanoic acid;

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1:1-mixture of (*rac.*)-(3*R**)-5-((*1R**, 5*S**)-6-[cyclopropyl-(3,5-dimethoxybenzyl)carbamoyl]-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-en-3-yl)-3-hydroxy-5-oxopentanoic acid and (*rac.*)-(3*S**)-5-((*1R**, 5*S**)-6-[cyclopropyl-(3,5-dimethoxybenzyl)carbamoyl]-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-en-3-yl)-3-hydroxy-5-oxopentanoic acid;

(*rac.*)-(1*R**, 5*S**)-3-acetyl-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid (2-chlorobenzyl)ethylamide;

20

(*rac.*)-(1*R**, 5*S**)-3-acetyl-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid cyclopropyl(3,5-dimethoxybenzyl)amide;

25 (*rac.*)-5-((1*R**, 5*S**)-6-(benzylcyclopropylcarbamoyl)-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-en-3-yl)-5-oxopentanoic acid;

30 (*rac.*)-5-((1*R**, 5*S**)-6-[(2-chlorobenzyl)ethylcarbamoyl]-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-en-3-yl)-5-oxopentanoic acid;

- (*rac.*)-5-((*IR**, *5S**)-6-[cyclopropyl-(2-fluorobenzyl)carbamoyl]-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-en-3-yl)-5-oxo-pentanoic acid;
- 5 (*rac.*)-5-((*IR**, *5S**)-6-[cyclopropyl-(3-trifluoromethylbenzyl)carbamoyl]-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-en-3-yl)-5-oxopentanoic acid;
- 10 (*rac.*)-5-((*IR**, *5S**)-6-[cyclopropyl(2-methylbenzyl)carbamoyl]-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-en-3-yl)-5-oxo-pentanoic acid;
- 15 (*rac.*)-5-((*IR**, *5S**)-6-{[2-(2-chlorophenyl)ethyl]cyclopropylcarbamoyl}-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-en-3-yl)-5-oxopentanoic acid;
- 20 (*rac.*)-5-((*IR**, *5S**)-6-[cyclopropyl(3,5-dimethoxybenzyl)carbamoyl]-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-en-3-yl)-5-oxopentanoic acid;
- 25 (*rac.*)-5-((*IR**, *5S**)-6-[cyclopropyl(2-*p*-tolylethyl)carbamoyl]-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-en-3-yl)-5-oxo-pentanoic acid;
- 30 (*rac.*)-5-((*IR**, *5S**)-6-[(2-chlorobenzyl)ethylcarbamoyl]-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-en-3-yl)-5-oxopentanoic acid methyl ester;
- (*rac.*)-5-((*IR**, *5S**)-6-[cyclopropyl-(2-fluorobenzyl)carbamoyl]-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-en-3-yl)-5-oxo-pentanoic acid methyl ester;

- (*rac.*)-5-((*1R*^{*}, *5S*^{*})-6-[cyclopropyl-(3,5-dimethoxybenzyl)carbamoyl]-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-en-3-yl)-5-oxopentanoic acid methyl ester;
- 5 (*rac.*)-5-((*1R*^{*}, *5S*^{*})-6-[(2-chlorobenzyl)ethylcarbamoyl]-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-en-3-yl)-2,2-dimethyl-5-oxopentanoic acid;
- 10 (*rac.*)-5-((*1R*^{*}, *5S*^{*})-6-[cyclopropyl-(2-fluorobenzyl)carbamoyl]-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-en-3-yl)-2,2-dimethyl-5-oxopentanoic acid;
- 15 (*rac.*)-5-((*1R*^{*}, *5S*^{*})-6-[cyclopropyl-(2-methylbenzyl)carbamoyl]-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-en-3-yl)-2,2-dimethyl-5-oxopentanoic acid;
- 20 (*rac.*)-5-((*1R*^{*}, *5S*^{*})-6-[cyclopropyl-(3,5-dimethoxybenzyl)carbamoyl]-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-en-3-yl)-2,2-dimethyl-5-oxopentanoic acid;
- 25 1:1-mixture of (*rac.*)-(2*R*^{*}, 3*S*^{*})-4-((*1R*^{*}, *5S*^{*})-6-[cyclopropyl-(2-methylbenzyl)carbamoyl]-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-en-3-yl)-2,3-dihydroxy-4-oxobutyric acid and (*rac.*)-(2*S*^{*}, 3*R*^{*})-4-((*1R*^{*}, *5S*^{*})-6-[cyclopropyl-(2-methylbenzyl)carbamoyl]-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-en-3-yl)-2,3-dihydroxy-4-oxobutyric acid;
- 30 (*rac.*)-5-((*1R*^{*}, *5S*^{*})-7-{4-[3-(2-bromo-5-fluorophenoxy)propyl]phenyl}-6-[cyclopropyl-(2-fluorobenzyl)carbamoyl]-3,9-diazabicyclo[3.3.1]non-6-en-3-yl)-5-oxopentanoic acid;

- (*rac.*)-5-{(*1R*^{*}, *5S*^{*})-7-{4-[3-(2-bromo-5-fluorophenoxy)propyl]phenyl}-6-[cyclopropyl-[2-(3-methylphenoxy)ethyl]carbamoyl]-3,9-diazabicyclo[3.3.1]-non-6-en-3-yl)-5-oxopentanoic acid;
- 5 (*rac.*)-5-[(*1R*^{*}, *5S*^{*})-7-{4-[3-(2-bromo-5-fluorophenoxy)propyl]phenyl}-6-(cyclopropylphenethylcarbamoyl)-3,9-diazabicyclo[3.3.1]non-6-en-3-yl]-5-oxopentanoic acid;
- 10 (*rac.*)-5-((*1R*^{*}, *5S*^{*})-7-{4-[3-(2-bromo-5-fluorophenoxy)propyl]phenyl}-6-{{2-(2-chlorophenyl)ethyl]cyclopropylcarbamoyl}-3,9-diazabicyclo[3.3.1]non-6-en-3-yl)-5-oxopentanoic acid;
- 15 (*rac.*)-5-((*1R*^{*}, *5S*^{*})-7-{4-[3-(2-bromo-5-fluorophenoxy)propyl]phenyl}-6-{{2-(2,3-difluorophenyl)ethyl]cyclopropylcarbamoyl}-3,9-diazabicyclo[3.3.1]non-6-en-3-yl)-5-oxopentanoic acid;
- 20 (*rac.*)-5-((*1R*^{*}, *5S*^{*})-7-{4-[3-(2-bromo-5-fluorophenoxy)propyl]phenyl}-6-{{2-(2-methylphenyl)ethyl]cyclopropylcarbamoyl}-3,9-diazabicyclo[3.3.1]non-6-en-3-yl)-5-oxopentanoic acid;
- 25 (*rac.*)-5-((*1R*^{*}, *5S*^{*})-7-{4-[3-(2-bromo-5-fluorophenoxy)propyl]phenyl}-6-{{2-(4-methylphenyl)ethyl]cyclopropylcarbamoyl}-3,9-diazabicyclo[3.3.1]non-6-en-3-yl)-5-oxopentanoic acid;
- 30 (*rac.*)-5-{(*1R*^{*}, *5S*^{*})-7-{4-[3-(2-bromo-5-fluorophenoxy)propyl]phenyl}-6-[cyclopropyl-(3,5-dimethoxybenzyl)carbamoyl]-3,9-diazabicyclo[3.3.1]non-6-en-3-yl}-5-oxopentanoic acid methyl ester;

- (*rac.*)-5-{(*1R*^{*}, *5S*^{*})-7-{4-[3-(2-bromo-5-fluorophenoxy)propyl]phenyl}-6-[cyclopropyl-(2-fluorobenzyl)carbamoyl]-3,9-diazabicyclo[3.3.1]non-6-en-3-yl}-2,2-dimethyl-5-oxopentanoic acid;
- 5 (*rac.*)-5-{(*1R*^{*}, *5S*^{*})-7-{4-[3-(2-bromo-5-fluorophenoxy)propyl]phenyl}-6-[cyclopropyl-(2-methylbenzyl)carbamoyl]-3,9-diazabicyclo[3.3.1]non-6-en-3-yl}-2,2-dimethyl-5-oxopentanoic acid;
- 10 (*rac.*)-5-[(*1R*^{*}, *5S*^{*})-7-{4-[3-(2-bromo-5-fluorophenoxy)propyl]phenyl}-6-(cyclopropylphenethylcarbamoyl)-3,9-diazabicyclo[3.3.1]non-6-en-3-yl]-2,2-dimethyl-5-oxopentanoic acid;
- 15 (*rac.*)-5-((*1R*^{*}, *5S*^{*})-7-{4-[3-(2-bromo-5-fluorophenoxy)propyl]phenyl}-6-{{[2-(4-methylphenyl)ethyl]cyclopropylcarbamoyl}-3,9-diazabicyclo[3.3.1]non-6-en-3-yl}-2,2-dimethyl-5-oxopentanoic acid;
- 20 (*rac.*)-5-{(*1R*^{*}, *5S*^{*})-7-{4-[3-(2-bromo-5-fluorophenoxy)propyl]phenyl}-6-[cyclopropyl-(3,5-dimethoxybenzyl)carbamoyl]-3,9-diazabicyclo[3.3.1]non-6-en-3-yl}-2,2-dimethyl-5-oxopentanoic acid;
- 25 1:1-mixture of (*rac.*)-(2*R*^{*}, 3*S*^{*})-4-{(*1R*^{*}, *5S*^{*})-7-{4-[3-(2-bromo-5-fluorophenoxy)propyl]phenyl}-6-[cyclopropyl-(3,5-dimethoxybenzyl)carbamoyl]-3,9-diazabicyclo[3.3.1]non-6-en-3-yl}-2,3-dihydroxy-4-oxobutyric acid and (*rac.*)-(2*S*^{*}, 3*R*^{*})-4-{(*1R*^{*}, *5S*^{*})-7-{4-[3-(2-bromo-5-fluorophenoxy)propyl]phenyl}-6-[cyclopropyl-(3,5-dimethoxybenzyl)carbamoyl]-3,9-diazabicyclo[3.3.1]non-6-en-3-yl}-2,3-dihydroxy-4-oxobutyric acid;
- 30 (*rac.*)-5-((*1R*^{*}, *5S*^{*})-6-[(2-chlorobenzyl)cyclopropylcarbamoyl]-7-{4-[2-(2,3,5-trimethylphenoxy)ethyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-en-3-yl)-5-oxo-pentanoic acid;

- (*rac.*.)-5-((*1R*^{*}, *5S*^{*})-6-[(3-trifluoromethylbenzyl)cyclopropylcarbamoyl]-7-{4-[2-(2,3,5-trimethylphenoxy)ethyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-en-3-yl)-5-oxopentanoic acid;
- 5 (*rac.*.)-5-((*1R*^{*}, *5S*^{*})-6-[(2-methylbenzyl)cyclopropylcarbamoyl]-7-{4-[2-(2,3,5-trimethylphenoxy)ethyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-en-3-yl)-5-oxopentanoic acid;
- 10 (*rac.*.)-5-((*1R*^{*}, *5S*^{*})-6-[(2-chlorobenzyl)ethylcarbamoyl]-7-{4-[2-(2,3,5-trimethylphenoxy)ethyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-en-3-yl)-2,2-dimethyl-5-oxopentanoic acid;
- 15 (*rac.*.)-5-((*1R*^{*}, *5S*^{*})-6-[cyclopropyl-(2-fluorobenzyl)carbamoyl]-7-{4-[2-(2,3,5-trimethylphenoxy)ethyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-en-3-yl)-2,2-dimethyl-5-oxopentanoic acid;
- 20 (*rac.*.)-5-((*1R*^{*}, *5S*^{*})-7-{4-[3-(2-bromo-5-fluorophenoxy)propyl]phenyl}-6-[(2-chlorobenzyl)cyclopropylcarbamoyl]-3,9-diazabicyclo[3.3.1]non-6-en-3-yl)-5-oxopentanoic acid;
- 25 (*rac.*.)-5-((*1R*^{*}, *5S*^{*})-7-{4-[3-(2-bromo-5-fluorophenoxy)propyl]phenyl}-6-[(2-chlorobenzyl)cyclopropylcarbamoyl]-3,9-diazabicyclo[3.3.1]non-6-en-3-yl)-2,2-dimethyl-5-oxopentanoic acid;
- 30 (*rac.*.)-5-((*1R*^{*}, *5S*^{*})-4-[3-(2-bromo-5-fluorophenoxy)propyl]phenyl)-3-(4-carbamoylbutyryl)-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid (2-chlorobenzyl)cyclopropylamide;
- (*rac.*.)-5-((*1R*^{*}, *5S*^{*})-6-(benzylcyclopropylcarbamoyl)-7-{4-[3-(2-bromo-5-fluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-en-3-yl)-5-oxopentanoic acid methyl ester;

- (*rac.*)-(1*R*^{*}, 5*S*^{*})-3-acetyl-7-{4-[2-(2-chloro-4,5-dimethylphenoxy)ethoxy]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid (2-chlorobenzyl)cyclopropylamide;
- 5 (*rac.*)-5-((1*R*^{*}, 5*S*^{*})-6-[(2-chlorobenzyl)cyclopropylcarbamoyl]-7-{4-[2-(2-chloro-4,5-dimethylphenoxy)ethoxy]phenyl}-3,9-diazabicyclo[3.3.1]non-6-en-3-yl)-5-oxopentanoic acid;
- 10 (*rac.*)-5-{(1*R*^{*}, 5*S*^{*})-7-{4-[2-(2-chloro-4,5-dimethylphenoxy)ethoxy]phenyl}-6-[cyclopropyl-(2-methylbenzyl)carbamoyl]-3,9-diazabicyclo[3.3.1]non-6-en-3-yl}-5-oxopentanoic acid;
- 15 (*rac.*)-5-{(1*R*^{*}, 5*S*^{*})-7-{4-[2-(2-chloro-4,5-dimethylphenoxy)ethoxy]phenyl}-6-[cyclopropyl-(3,5-dimethoxybenzyl)carbamoyl]-3,9-diazabicyclo[3.3.1]non-6-en-3-yl}-5-oxopentanoic acid;
- 20 (*rac.*)-5-((1*R*^{*}, 5*S*^{*})-6-[(2-chlorobenzyl)cyclopropylcarbamoyl]-7-{4-[2-(2-chloro-4,5-dimethylphenoxy)ethoxy]phenyl}-3,9-diazabicyclo[3.3.1]non-6-en-3-yl)-2,2-dimethyl-5-oxopentanoic acid;
- 25 (*rac.*)-5-((1*R*^{*}, 5*S*^{*})-6-[(2-chlorobenzyl)ethylcarbamoyl]-7-{4-[2-(2-chloro-4,5-dimethylphenoxy)ethoxy]phenyl}-3,9-diazabicyclo[3.3.1]non-6-en-3-yl)-2,2-dimethyl-5-oxopentanoic acid;
- 30 (*rac.*)-5-{(1*R*^{*}, 5*S*^{*})-7-{4-[2-(2-chloro-4,5-dimethylphenoxy)ethoxy]phenyl}-6-[cyclopropyl-(2-fluorobenzyl)carbamoyl]-3,9-diazabicyclo[3.3.1]non-6-en-3-yl}-2,2-dimethyl-5-oxopentanoic acid;
- 35 (*rac.*)-5-((1*R*^{*}, 5*S*^{*})-6-[cyclopropyl-(3,5-dimethoxybenzyl)carbamoyl]-7-{4-[2-(2,3,5-trimethylphenoxy)ethyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-en-3-yl)-5-oxopentanoic acid;

- (*rac.*)-5-((*1R*^{*}, *5S*^{*})-6-[cyclopropyl-(2-*p*-tolylethyl)carbamoyl]-7-{4-[2-(2,3,5-trimethylphenoxy)ethyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-en-3-yl)-5-oxopentanoic acid;
- 5 (*rac.*)-5-((*1R*^{*}, *5S*^{*})-7-{4-[2-(2-chloro-4,5-dimethylphenoxy)ethoxy]phenyl}-6-{cyclopropyl-[2-(2-hydroxyethyl)benzyl]carbamoyl}-3,9-diazabicyclo[3.3.1]non-6-en-3-yl)-5-oxopentanoic acid;
- 10 (*rac.*)-5-((*1R*^{*}, *5S*^{*})-6-(cyclopropylphenethylcarbamoyl)-7-{4-[2-(2,3,5-trimethylphenoxy)ethyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-en-3-yl)-2,2-dimethyl-5-oxopentanoic acid;
- 15 (*rac.*)-(*1R*^{*}, *5S*^{*})-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid (2-chlorobenzyl)ethylamide;
- (*rac.*)-(*1R*^{*}, *5S*^{*})-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-3,6-dicarboxylic acid 6-[(2-chlorobenzyl)cyclopropylamide] 3-dimethylamide;
- 20 (*rac.*)-(*1R*^{*}, *5S*^{*})-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-3,6-dicarboxylic acid 6-[(2-chlorobenzyl)cyclopropylamide] 3-diethylamide;
- 25 (*rac.*)-(*1R*^{*}, *5S*^{*})-6-[(2-chlorobenzyl)cyclopropylcarbamoyl]-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-3-carboxylic acid methyl ester;
- 30 (*rac.*)-(*1R*^{*}, *5S*^{*})-6-[(2-chlorobenzyl)cyclopropylcarbamoyl]-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-3-carboxylic acid ethyl ester;

(*rac.*)-(1*R*^{*}, 5*S*^{*})-3-methanesulfonyl-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]-phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid (2-chlorobenzyl)-cyclopropylamide;

5 (*rac.*)-(1*R*^{*}, 5*S*^{*})-3-ethanesulfonyl-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]-phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid (2-chlorobenzyl)-cyclopropylamide;

10 (*rac.*)-5-((1*R*^{*}, 5*S*^{*})-6-[(2-chlorobenzyl)cyclopropylcarbamoyl]-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-en-3-yl)-5-oxo-pentanoic acid ethyl ester;

15 (*rac.*)-4-((1*R*^{*}, 5*S*^{*})-6-[(2-chlorobenzyl)cyclopropylcarbamoyl]-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-en-3-yl)-4-oxo-butrylic acid;

20 (*rac.*)-3-[((1*R*^{*}, 5*S*^{*})-6-[(2-chlorobenzyl)cyclopropylcarbamoyl]-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-3-carbonyl)-amino]propionic acid ethyl ester;

25 (*rac.*)-4-[((1*R*^{*}, 5*S*^{*})-6-[(2-chlorobenzyl)cyclopropylcarbamoyl]-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-3-carbonyl)-amino]butyric acid ethyl ester;

30 (*rac.*)-(1*R*^{*}, 5*S*^{*})-3-(3-carbamoylpropionyl)-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid (2-chlorobenzyl)cyclopropylamide;

35 (*rac.*)-(1*R*^{*}, 5*S*^{*})-3-(2-hydroxyacetyl)-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]-phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid (2-chlorobenzyl)-cyclopropylamide;

(*1S*, *5R*)-3-((*3R*)-3-hydroxybutyryl)-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid (2-chlorobenzyl)cyclopropylamide;

- 5 1:1-mixture of (*rac.*)-(1*R*^{*}, 5*S*^{*})-3-((1*R*^{*}, 2*S*^{*})-2-hydroxycyclopentanecarbonyl)-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid (2-chlorobenzyl)cyclopropylamide and (*rac.*)-(1*R*^{*}, 5*S*^{*})-3-((1*S*^{*}, 2*R*^{*})-2-hydroxycyclopentanecarbonyl)-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid (2-chlorobenzyl)cyclopropylamide;

10 (15) (*rac.*)-(1*R*^{*}, 5*S*^{*})-9-acetyl-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid (2-chlorobenzyl)cyclopropylamide;

15 (16) (*rac.*)-5-((1*R*^{*}, 5*S*^{*})-6-[(2-chlorobenzyl)cyclopropylcarbamoyl]-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-en-9-yl)-5-oxo-pentanoic acid;

20 (17) (20) (*rac.*)-5-((1*R*^{*}, 5*S*^{*})-6-[(2-chlorobenzyl)cyclopropylcarbamoyl]-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-en-9-yl)-5-oxo-pentanoic acid ethyl ester;

25 (18) (25) (*rac.*)-5-((1*R*^{*}, 5*S*^{*})-6-[(2-chlorobenzyl)cyclopropylcarbamoyl]-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-en-9-yl)-5-oxo-pentanoic acid methyl ester;

30 (19) 1:1-mixture of (*rac.*)-(3*R*^{*})-5-((1*R*^{*}, 5*S*^{*})-6-[(2-chlorobenzyl)cyclopropylcarbamoyl]-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-en-9-yl)-3-hydroxy-5-oxo-pentanoic acid and (*rac.*)-(3*S*^{*})-5-((1*R*^{*}, 5*S*^{*})-6-[(2-chlorobenzyl)cyclopropylcarbamoyl]-7-{4-[3-(2,3,6-trifluoro-

phenoxy)propyl]phenyl}-3,9-diazabicyclo-[3.3.1]non-6-en-9-yl)-3-hydroxy-5-oxopentanoic acid;

5 (rac.)-5-((1*R*^{*}, 5*S*^{*})-6-[(2-chlorobenzyl)cyclopropylcarbamoyl]-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-en-9-yl)-2,2-dimethyl-5-oxopentanoic acid;

10 (rac.)-4-((1*R*^{*}, 5*S*^{*})-6-[(2-chlorobenzyl)cyclopropylcarbamoyl]-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-en-9-yl)-4-oxo-15 butyric acid;

15 (rac.)-(1*R*^{*}, 5*S*^{*})-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6,9-dicarboxylic acid 6-[(2-chlorobenzyl)cyclopropylamide] 9-dimethylamide;

20 (rac.)-(1*R*^{*}, 5*S*^{*})-6-[(2-chlorobenzyl)cyclopropylcarbamoyl]-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-9-carboxylic acid methyl ester;

25 (rac.)-(1*R*^{*}, 5*S*^{*})-6-[(2-chlorobenzyl)cyclopropylcarbamoyl]-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-9-carboxylic acid ethyl ester;

30 (rac.)-3-[(1*R*^{*}, 5*S*^{*})-(6-[(2-chlorobenzyl)cyclopropylcarbamoyl]-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-9-carbonyl)-amino]propionic acid ethyl ester;

(rac.)-4-[((1*R*^{*}, 5*S*^{*})-6-[(2-chlorobenzyl)cyclopropylcarbamoyl]-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-9-carbonyl)-amino]butyric acid ethyl ester;

- (*rac.*)-(1*R*^{*}, 5*S*^{*})-3-formyl-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid (2-chlorobenzyl)cyclopropylamide;
- 5 (*rac.*)-3-[((1*R*^{*}, 5*S*^{*})-6-[(2-chlorobenzyl)cyclopropylcarbamoyl]-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-3-carbonyl)-amino]propionic acid;
- 10 (*rac.*)-3-[((1*R*^{*}, 5*S*^{*})-6-[(2-chlorobenzyl)cyclopropylcarbamoyl]-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-9-carbonyl)-amino]propionic acid;
- 15 (*rac.*)-4-[((1*R*^{*}, 5*S*^{*})-6-[(2-chlorobenzyl)cyclopropylcarbamoyl]-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-9-carbonyl)-amino]butyric acid;
- 20 (*rac.*)-(1*R*^{*}, 5*S*^{*})-3-acetyl-7-{4-[2-(2,3,6-trifluorophenoxy)ethyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid (2-chlorobenzyl)cyclopropylamide;
- 25 (*rac.*)-(1*R*^{*}, 5*S*^{*})-9-(4-carbamoylbutyryl)-7-{4-[2-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid (2-chlorobenzyl)cyclopropylamide;
- 30 (*rac.*)-(1*R*^{*}, 5*S*^{*})-9-(3-carbamoylpropionyl)-7-{4-[2-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid (2-chlorobenzyl)cyclopropylamide;
- 30 (*rac.*)-(1*R*^{*}, 5*S*^{*})-9-(2-hydroxyacetyl)-7-{4-[2-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid (2-chlorobenzyl)cyclopropylamide;

- (*1R*, *5S*)-9-((*3S*)-3-hydroxybutyryl)-7-{4-[2-(2,3,6-trifluorophenoxy)propyl]-phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid (2-chlorobenzyl)-cyclopropylamide;
- 5 (*rac.*)-(*1R**, *5S**)-9-methanesulfonyl-7-{4-[2-(2,3,6-trifluorophenoxy)propyl]-phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid (2-chlorobenzyl)-cyclopropylamide;
- (*rac.*)-(*1R**, *5S**)-9-ethanesulfonyl-7-{4-[2-(2,3,6-trifluorophenoxy)propyl]-phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid (2-chlorobenzyl)-cyclopropylamide;
- 10 (*rac.*)-(*1R**, *5S**)-3-acetyl-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid cyclopropyl-(2-fluoro-5-methoxybenzyl)amide;
- 15 (*rac.*)-(*1R**, *5S**)-3-acetyl-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid cyclopropyl-(3-methoxybenzyl)amide;
- 20 (*rac.*)-(*1R**, *5S**)-3-acetyl-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid cyclopropyl-(3,4-dimethoxybenzyl)amide;
- 25 (*rac.*)-(*1R**, *5S**)-3-acetyl-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid (2-chloro-3-trifluoromethylbenzyl)cyclopropylamide;
- (*rac.*)-(*1R**, *5S**)-3-acetyl-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid benzo[1,3]dioxol-5-ylmethyl-cyclopropylamide;
- 30

- (*rac.*)-(1*R*^{*}, 5*S*^{*})-3-acetyl-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid (2-chloro-6-fluorobenzyl)cyclopropylamide;
- 5 (*rac.*)-(1*R*^{*}, 5*S*^{*})-3-acetyl-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid (2-bromobenzyl)cyclopropylamide;
- 10 (*rac.*)-(1*R*^{*}, 5*S*^{*})-3-acetyl-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid cyclopropyl-(2,3-dimethylbenzyl)amide;
- 15 (*rac.*)-(1*R*^{*}, 5*S*^{*})-3-acetyl-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid cyclopropyl-(3,5-difluorobenzyl)amide;
- 20 (*rac.*)-(1*R*^{*}, 5*S*^{*})-3-acetyl-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid cyclopropyl-(2,3-dichlorobenzyl)amide;
- 25 (*rac.*)-(1*R*^{*}, 5*S*^{*})-3-acetyl-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid cyclopropyl-(3-trifluoromethoxybenzyl)amide;
- 30 (*rac.*)-(1*R*^{*}, 5*S*^{*})-3-acetyl-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid cyclopropyl-(3-methylbenzyl)amide;
- (*rac.*)-(1*R*^{*}, 5*S*^{*})-3-acetyl-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid (3-chlorobenzyl)cyclopropylamide;

- (*rac.*)-(1*R*^{*}, 5*S*^{*})-3-(5-morpholin-4-yl-5-oxopentanoyl)-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid (2-chlorobenzyl)cyclopropylamide;
- 5 (*rac.*)-(1*R*^{*}, 5*S*^{*})-3-(2-tetrazol-1-ylacetyl)-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid (2-chlorobenzyl)cyclopropylamide;
- 10 (*rac.*)-(1*R*^{*}, 5*S*^{*})-3-(5-oxo-5-piperazin-1-ylpentanoyl)-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid (2-chlorobenzyl)cyclopropylamide;
- 15 1:1-mixture of (1*R*, 5*S*)-3-((2*S*)-2-amino-3-hydroxypropionyl)-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid (2-chlorobenzyl)cyclopropylamide and (1*S*, 5*R*)-3-((2*S*)-2-amino-3-hydroxypropionyl)-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid (2-chlorobenzyl)cyclopropylamide;
- 20 1:1-mixture of (1*R*, 5*S*)-3-((2*S*)-2-aminopropionyl)-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid (2-chlorobenzyl)cyclopropylamide and (1*S*, 5*R*)-3-((2*S*)-2-aminopropionyl)-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid (2-chlorobenzyl)cyclopropylamide;
- 25 (*rac.*)-(1*R*^{*}, 5*S*^{*})-3-acetyl-7-{4-[2-(2,6-dichloro-4-methylphenoxy)ethoxy]-phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid cyclopropyl-(2,3-dichlorobenzyl)amide;
- 30 (*rac.*)-(1*R*^{*}, 5*S*^{*})-3-acetyl-7-{4-[2-(2,6-dichloro-4-methylphenoxy)ethoxy]-phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid cyclopropyl-(2,3-dimethylbenzyl)amide;

(*rac.*)-(1*R*^{*}, 5*S*^{*})-3-acetyl-7-{4-[2-(2,6-dichloro-4-methylphenoxy)ethoxy]-phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid (2-chloro-3-trifluoromethyl-benzyl)cyclopropylamide;

5 (*rac.*)-(1*R*^{*}, 5*S*^{*})-3-acetyl-7-{4-[2-(2,6-dichloro-4-methylphenoxy)ethoxy]-phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid (2-bromobenzyl)-cyclopropylamide;

10 (*rac.*)-(1*R*^{*}, 5*S*^{*})-3-acetyl-7-{4-[2-(2,6-dichloro-4-methylphenoxy)ethoxy]-phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid (2-chlorobenzyl)-cyclopropylamide;

15 (*rac.*)-(1*R*^{*}, 5*S*^{*})-3-acetyl-7-{4-[2-(2,6-dichloro-4-methylphenoxy)ethoxy]-phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid (2-chlorobenzyl)-ethylamide;

20 (*rac.*)-(1*R*^{*}, 5*S*^{*})-3-acetyl-7-{4-[2-(2,6-dichloro-4-methylphenoxy)ethoxy]-phenyl}-3,9-diaza-bicyclo[3.3.1]non-6-ene-6-carboxylic acid (6-chlorobenzo-[1,3]dioxol-5-ylmethyl)cyclopropylamide;

25 (*rac.*)-(1*R*^{*}, 5*S*^{*})-3-acetyl-7-{4-[2-(2,6-dichloro-4-methylphenoxy)ethoxy]-phenyl}-3,9-diaza-bicyclo[3.3.1]non-6-ene-6-carboxylic acid cyclopropyl-(3,5-dimethoxybenzyl)amide;

30 (*rac.*)-(1*R*^{*}, 5*S*^{*})-3-acetyl-7-{4-[2-(2,6-dichloro-4-methylphenoxy)ethoxy]-phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid cyclopropyl-(3-methoxybenzyl)amide;

35 (*rac.*)-(1*R*^{*}, 5*S*^{*})-3-acetyl-7-{4-[2-(2-chloro-5-fluorophenoxy)ethoxy]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid cyclopropyl-(2,3-dichloro-benzyl)amide;

- (*rac.*)-(1*R*^{*}, 5*S*^{*})-3-acetyl-7-{4-[2-(2-chloro-5-fluorophenoxy)ethoxy]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid (2-chloro-3-trifluoromethylbenzyl)cyclopropylamide;
- 5 (*rac.*)-(1*R*^{*}, 5*S*^{*})-3-acetyl-7-{4-[2-(2,6-dichloro-4-methylphenoxy)ethoxy]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid (3-chlorobenzyl)cyclopropylamide;
- 10 (*rac.*)-(1*R*^{*}, 5*S*^{*})-3-acetyl-7-{4-[2-(2,6-dichloro-4-methylphenoxy)ethoxy]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid cyclopropyl-(3-methylbenzyl)amide;
- 15 (*rac.*)-(1*R*^{*}, 5*S*^{*})-3-acetyl-7-{4-[2-(2,6-dichloro-4-methylphenoxy)ethoxy]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid cyclopropyl-(2-fluoro-5-methoxybenzyl)amide;
- 20 (*rac.*)-(1*R*^{*}, 5*S*^{*})-3-acetyl-7-{4-[2-(2-chloro-4,5-dimethylphenoxy)ethoxy]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid cyclopropyl-(2,3-dichlorobenzyl)amide;
- 25 (*rac.*)-(1*R*^{*}, 5*S*^{*})-3-acetyl-7-{4-[3-(2-chloro-3,6-difluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid (2-chlorobenzyl)cyclopropylamide;
- 30 (*rac.*)-(1*R*^{*}, 5*S*^{*})-3-acetyl-7-{4-[2-(2-chloro-4,5-dimethylphenoxy)ethoxy]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid (2-chloro-3-trifluoromethylbenzyl)cyclopropylamide;
- (*rac.*)-(1*R*^{*}, 5*S*^{*})-3-acetyl-7-{4-[2-(2,6-dichloro-4-methylphenoxy)ethoxy]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid cyclopropyl-(3,4-dimethoxybenzyl)amide;

- (*rac.*)-(1*R*^{*}, 5*S*^{*})-3-acetyl-7-{4-[2-(2,6-dichloro-4-methylphenoxy)ethoxy]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid cyclopropyl-(3-trifluoromethoxybenzyl)amide;
- 5 (*rac.*)-(1*R*^{*}, 5*S*^{*})-3-acetyl-7-{4-[2-(2,4,5-trichlorophenoxy)ethoxy]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid (2-chloro-3-trifluoromethylbenzyl)cyclopropylamide;
- 10 (*rac.*)-(1*R*^{*}, 5*S*^{*})-3-acetyl-7-{4-[2-(2,4,5-trichlorophenoxy)ethoxy]phenyl}-3,9-diazabicyclo-[3.3.1]non-6-ene-6-carboxylic acid cyclopropyl-(2,3-dichlorobenzyl)amide;
- 15 (*rac.*)-(1*R*^{*}, 5*S*^{*})-3-acetyl-7-{4-[2-(2,3-dichlorophenoxy)ethoxy]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid cyclopropyl-(2,3-dimethylbenzyl)amide;
- 20 (*rac.*)-(1*R*^{*}, 5*S*^{*})-3-acetyl-7-{4-[2-(2,4,5-trichlorophenoxy)ethoxy]phenyl}-3,9-diazabicyclo-[3.3.1]non-6-ene-6-carboxylic acid cyclopropyl-(2,3-dimethylbenzyl)amide;
- 25 (*rac.*)-(1*R*^{*}, 5*S*^{*})-3-acetyl-7-{4-[2-(2-chloro-4,5-dimethylphenoxy)ethoxy]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid cyclopropyl-(2,3-dimethylbenzyl)amide;
- 30 (*rac.*)-(1*R*^{*}, 5*S*^{*})-3-acetyl-7-{4-[2-(2-bromo-5-fluorophenoxy)ethoxy]phenyl}-3,9-diaza-bicyclo[3.3.1]non-6-ene-6-carboxylic acid (2-bromobenzyl)-cyclopropylamide;

(*rac.*)-(1*R*^{*}, 5*S*^{*})-3-acetyl-7-{4-[2-(2,3-dichlorophenoxy)ethoxy]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid cyclopropyl-(2,3-dichlorobenzyl)amide;

5 (*rac.*)-(1*R*^{*}, 5*S*^{*})-3-acetyl-7-{4-[2-(2-chloro-5-fluorophenoxy)ethoxy]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid cyclopropyl-(2,3-dimethylbenzyl)amide;

10 (*rac.*)-(1*R*, 5*S*)-7-{4-[3-(2,3,6-trifluorophenoxy)propylphenyl]-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid (2-chlorobenzyl)cyclopropylamide;

15 (*rac.*)-(1*S*, 5*R*)-7-{4-[3-(2,3,6-trifluorophenoxy)propylphenyl]-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid (2-chlorobenzyl)cyclopropylamide.

15 The compounds of general formula I and their pharmaceutically acceptable salts may be used as therapeutics e.g. in form of pharmaceutical compositions. They may especially be used in the treatment and/or prophylaxis of cardiovascular and renal diseases. Examples of such diseases are hypertension, coronary diseases, cardiac insufficiency, renal insufficiency, renal and myocardial ischemia, and 20 renal failure. They can also be used to prevent restenosis after balloon or stent angioplasty, to treat erectile dysfunction, glomerulonephritis, renal colic, and glaucoma. Furthermore, they can be used in the therapy and the prophylaxis of diabetic complications, complications of vascular or cardiac surgery or after organ transplantation, complications of cyclosporin treatment, as well as other diseases 25 presently known to be related to the RAS.

In another embodiment, the invention relates to a method for the treatment and/or prophylaxis of diseases which are related to the RAS such as hypertension, coronary diseases, cardiac insufficiency, renal insufficiency, renal and myocardial 30 ischemia, and renal failure, which method comprises administrating a compound as defined above to a human being or animal.

The invention further relates to the use of compounds of general formula I as defined above for the treatment and/or prophylaxis of diseases which are associated with the RAS such as hypertension, coronary diseases, cardiac insufficiency, renal insufficiency, renal and myocardial ischemia, and renal failure.

In addition, the invention relates to the use of compounds as defined above for the preparation of medicaments for the treatment and/or prophylaxis of diseases which are associated with the RAS such as hypertension, coronary diseases, cardiac insufficiency, renal insufficiency, renal and myocardial ischemia, and renal failure.

The compounds of formula I may also be used in combination with one or more other therapeutically useful substances e. g. with other renin inhibitors, with ACE-inhibitors, with angiotensin-receptor antagonists, with diuretics, with calcium channel blockers, with endothelin receptors antagonists or with other drugs beneficial for the prevention or the treatment of cardiovascular events or renal insufficiency.

All forms of prodrugs leading to an active component comprised by general formula I above are included in the present invention.

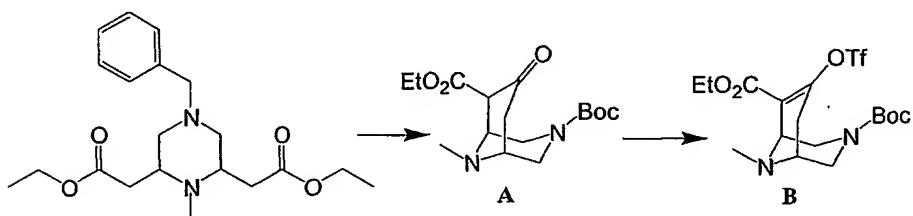
The compounds of general formula I can be manufactured by the methods outlined below, by the methods described in the examples or by analogous methods.

Preparation of the precursors:

Precursors are compounds which were prepared as key intermediates and/or building blocks and which were suitable for further transformations in parallel chemistry.

Bicyclonanonone **A** was prepared from (4-benzyl-6-ethoxycarbonylmethyl-1-methylpiperazin-2-yl)acetic acid ethyl ester (Patent Application WO92/05174) as described in Scheme 1. Derivative **A** might also be present as enol form. In order to allow a coupling at the 7-position of bicyclononane **A** with aryl bromides, the 5 vinyl triflate derivative **B** was prepared.

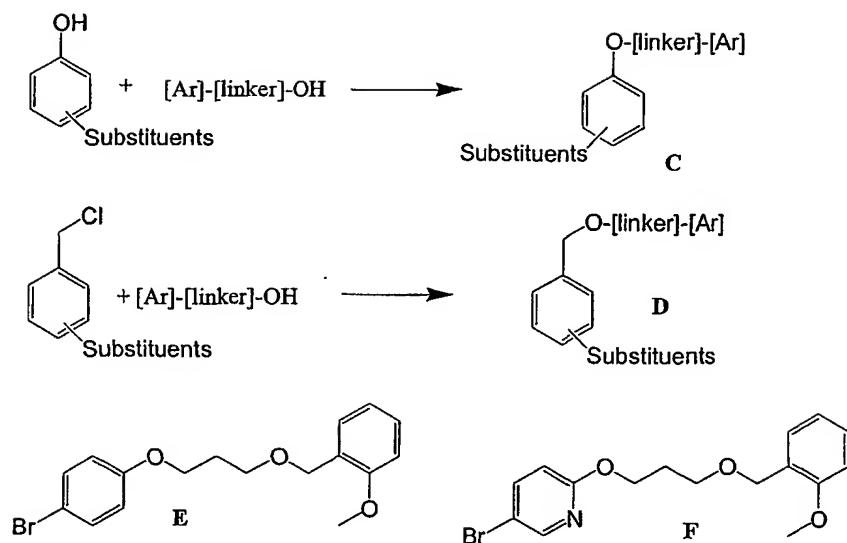
Scheme 1



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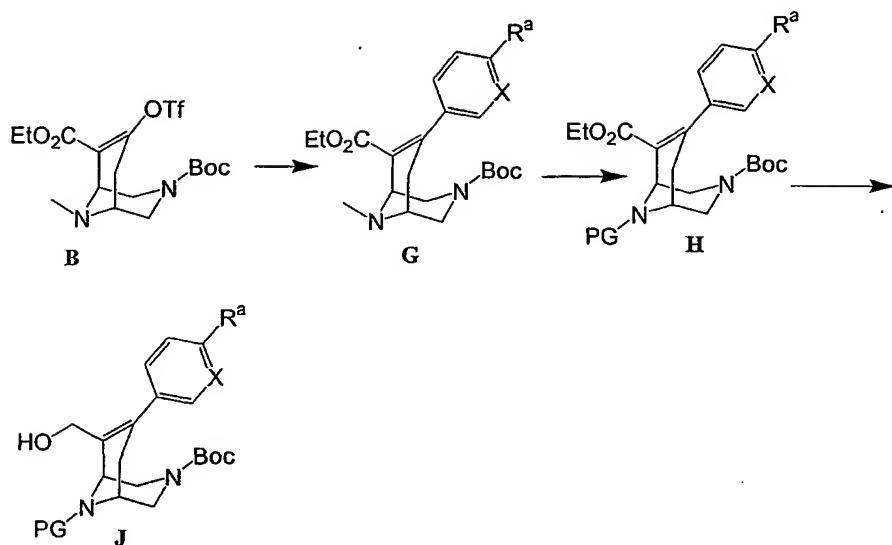
The bromoaryl components can be prepared as described in Scheme 2. A *Mitsunobu* coupling (\rightarrow compounds of type **C**) or the alkylation of an alcohol with a benzylic chloride (or bromide, \rightarrow compounds of type **D**) are often the most convenient methods. Derivatives **E** and **F** were prepared in one step from 1-(3-chloropropoxymethyl)-2-methoxybenzene (Vieira E. et al., *Bioorg. Med. Chem. Letters*, 1999, 9, 1397) or 3-(5-bromopyridin-2-yloxy)propan-1-ol (Patent Application WO 98/39328) according to these methods. Other methods for the preparation of ethers or thioethers, like a *Williamson* synthesis, might be used as well (see e.g. March, J, "Advanced Organic Chemistry", 3rd ed., John Wiley and sons, 1985).

Scheme 2



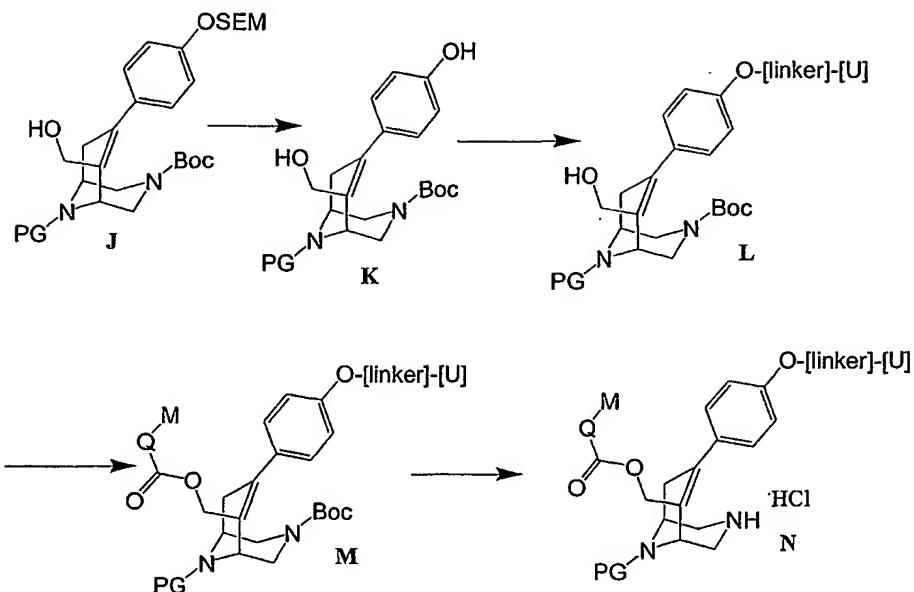
5 As shown in Scheme 3, these bromoaryl derivatives might then be coupled to triflate **B** in the presence of $Pd(PPh_3)_4$ or a related $Pd(0)$ -complex in order to obtain bicyclononenes **G** (for details see the corresponding examples). Protective group manipulation would lead to the bicyclononenes **H**, which can finally be reduced to the alcohol derivatives **J**.

Scheme 3



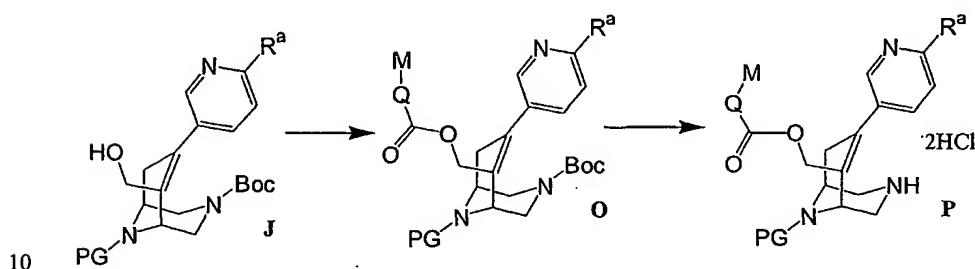
- 5 If R^a is O-SEM and X is CH, in compounds of type J, the SEM-protecting group
can be cleaved under slightly acidic conditions, while the Boc-protecting group
remains untouched as illustrated in Scheme 4. The phenolic moiety of
bicyclononene K might then be alkylated to bicyclononene L. This alkohol
intermediate would be transformed into the ester M, and the Boc-protecting group
10 can finally be cleaved to yield precursor N.

Scheme 4



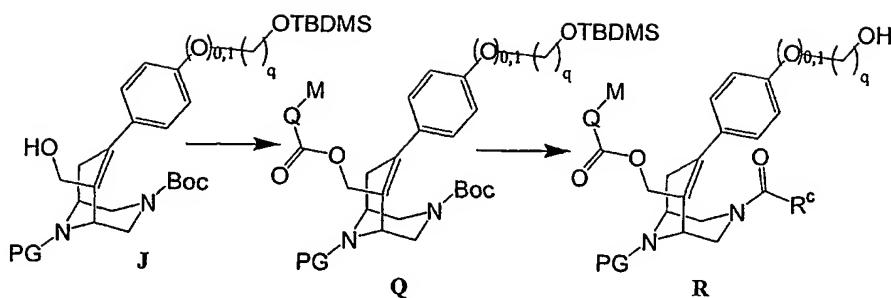
5 If X is N (Scheme 3), bicyclononenes of type J can be esterified to bicyclononenes O that can be deprotected to precursors P (Scheme 5).

Scheme 5



If, in alcohols J, X is CH (Scheme 6) and, for instance, R^a is O(CH₂)_qOTBDMS or (CH₂)_qOTBDMS (Scheme 3), an esterification might lead to bicyclononenes Q (Scheme 6). Under acidic conditions, both the Boc- and the TBDMS-groups would be cleaved and the secondary amine might be acylated, sulfonated, or alkylated to yield precursors R.

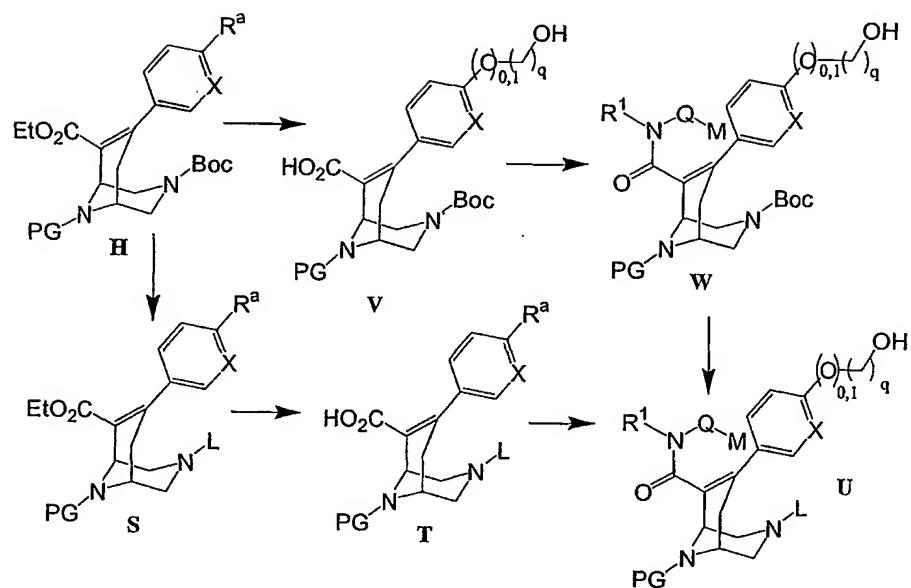
Scheme 6



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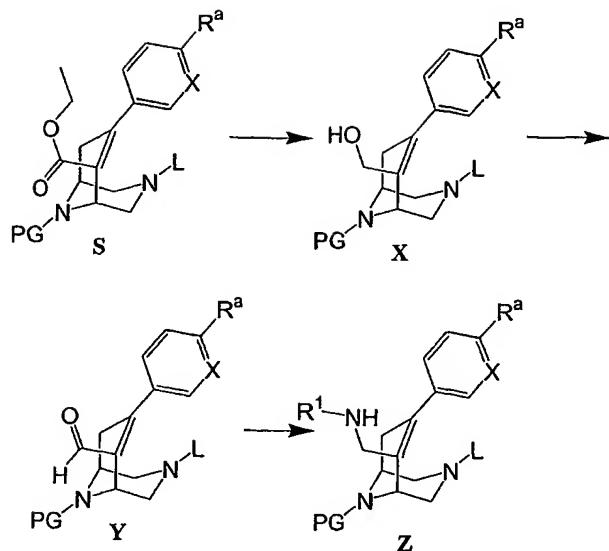
As illustrated in Scheme 7, the Boc-protecting group of bicyclononenes **H** might be cleaved and the resulting secondary amine acylated, sulfonated, or *N*-alkylated to bicyclononenes **S**. Saponification of the ester would lead to precursors **T**. If, for instance, R^a is $O(CH_2)_nOTBDMS$ or $(CH_2)_nOTBDMS$, the silyl ether might be 10 cleaved simultaneously during the cleavage of the Boc-protecting group or during the saponification. The acid might be transformed into an amide to lead to precursors **U**. Alternatively, amides **U** can be prepared from bicyclononenes **H** through the acids **V**, with simultaneous cleavage of the silyl ether. Reaction of the acids **V** with amines might give the amides **W** that can be transformed into 15 precursors **U**. The amines can be prepared according to the literature or as described in the experimental part.

Scheme 7



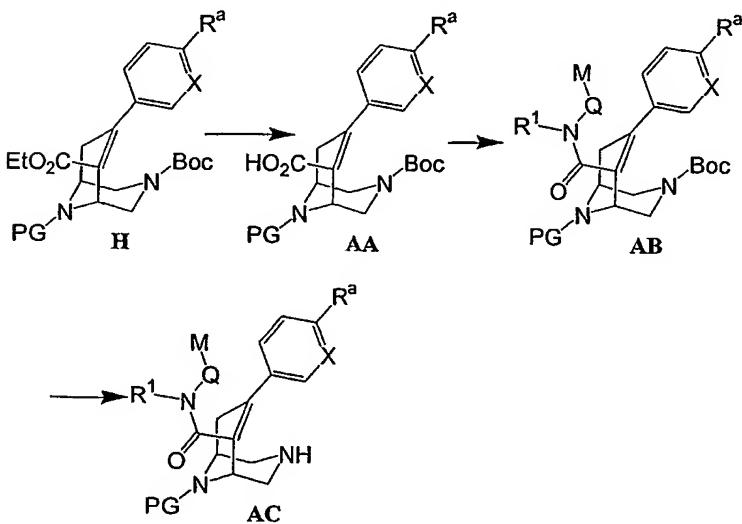
- 5 As illustrated in Scheme 8, the bicyclononenes **S** might be reduced to the corresponding alcohols **X**. The alcohol derivative **X** might then be oxidised to the aldehydes **Y**, which might transformed to the precursors **Z** by reductive amination.

Scheme 8



5 As shown in Scheme 9, a compound of type **H** might also be saponified to a compound of type **AA**. After amide coupling to a derivative of type **AB**, removal of the Boc-group would yield a precursor of type **AC**.

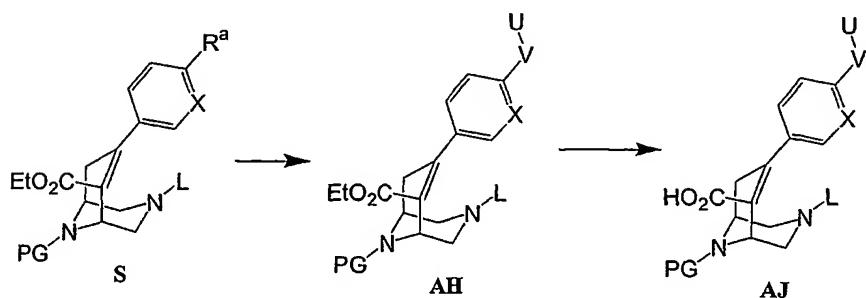
Scheme 9



As illustrated in Scheme 10, a compound of type **S** might also be transformed into a compound of type **AH** that in turn can be saponified to a precursor of type **AJ**.

Scheme 10

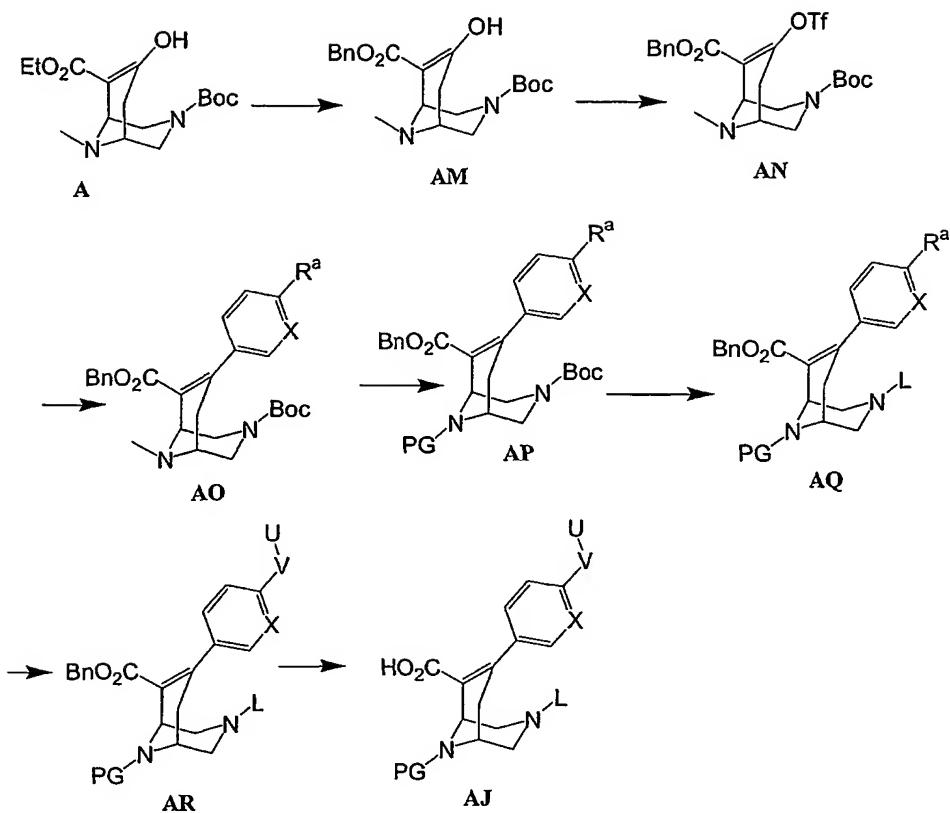
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Alternatively precursors of type **AJ** might be prepared from bicyclonanonane **A**, but using the benzyl ester instead of the ethyl ester, as illustrated in Scheme 11.

- 10 After a transesterification to compound **AM** similar reactions as described here above would lead to compounds **AN**, **AO**, **AP**, **AQ**, **AR** and finally **AJ**.

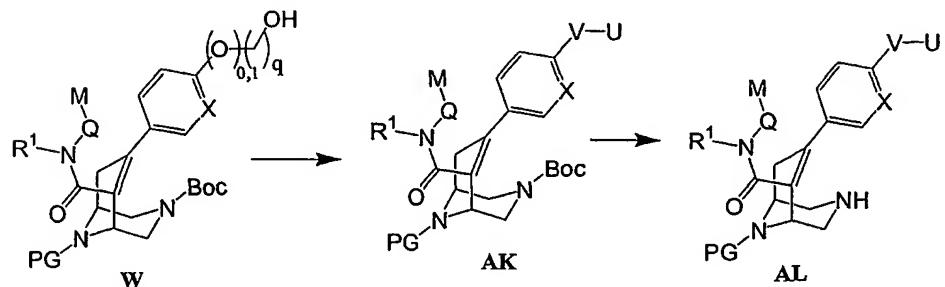
Scheme 11



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As shown in Scheme 12 a precursor of type AL might be prepared as well in two steps from a compound of type W.

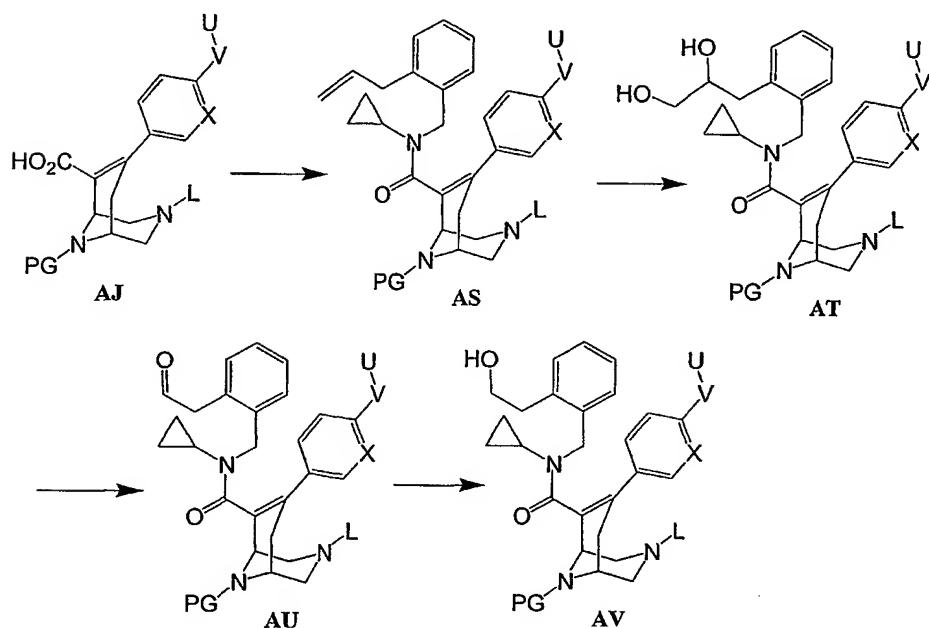
Scheme 12



Sometimes it might be necessary to transform a substituent further after attaching it on the bicyclic template. For instance a compound of type AS, obtained by amide coupling from a compound of type AJ, might be transformed into a precursor of type AT, AU, or AV, as illustrated in Scheme 13.

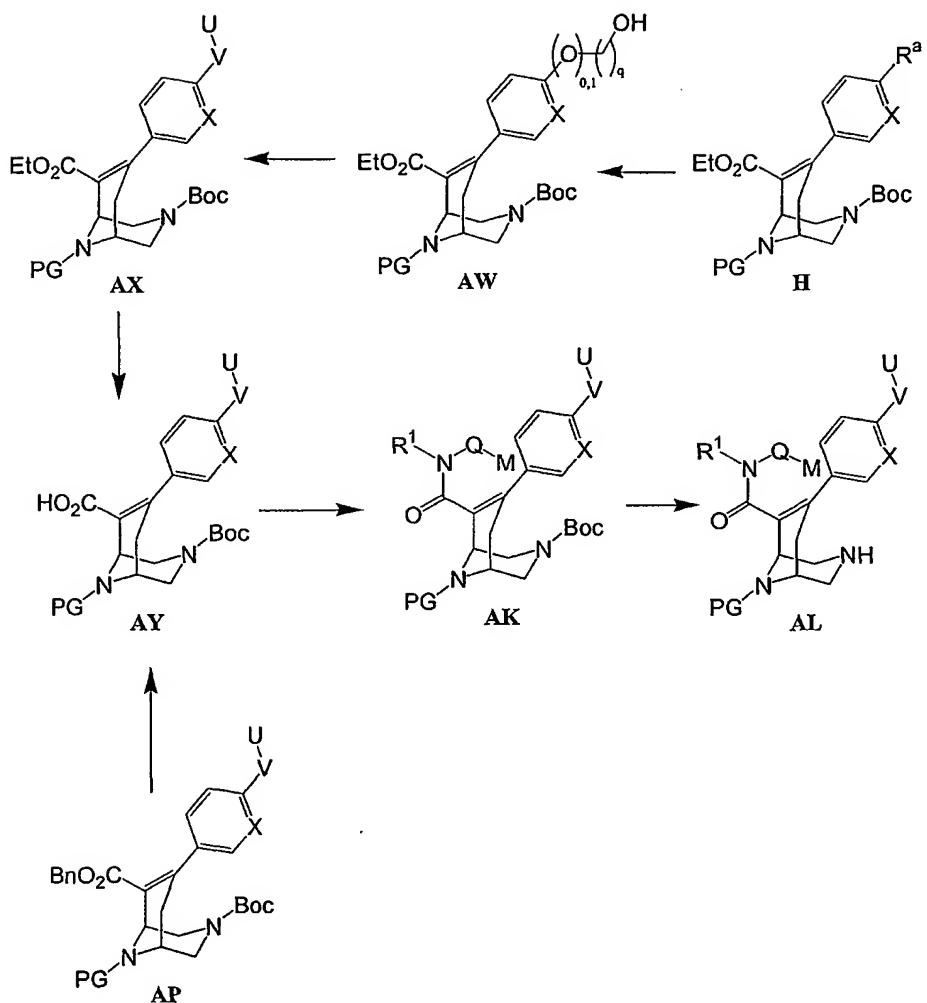
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Scheme 13



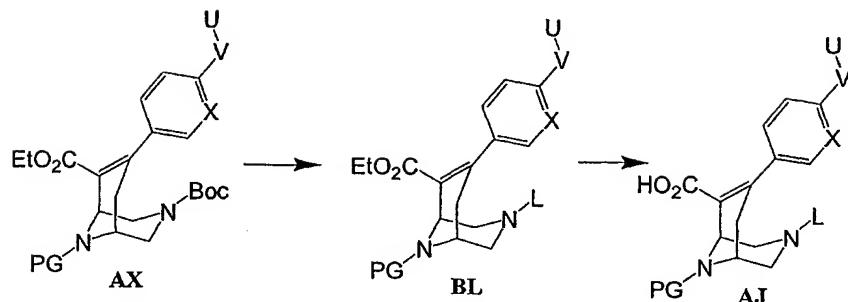
- 10 As illustrated in Scheme 14 compounds of type H might be deprotected into compounds of type AW. This type of compounds might be then transformed into compounds of type AX and finally into compounds of type AY. Compounds of type AY might also be prepared from compounds of type AP. Compounds of type AY might be transformed into compounds of type AK. Compounds of type AK might be finally transformed into precursors of type AL.
- 15

Scheme 14



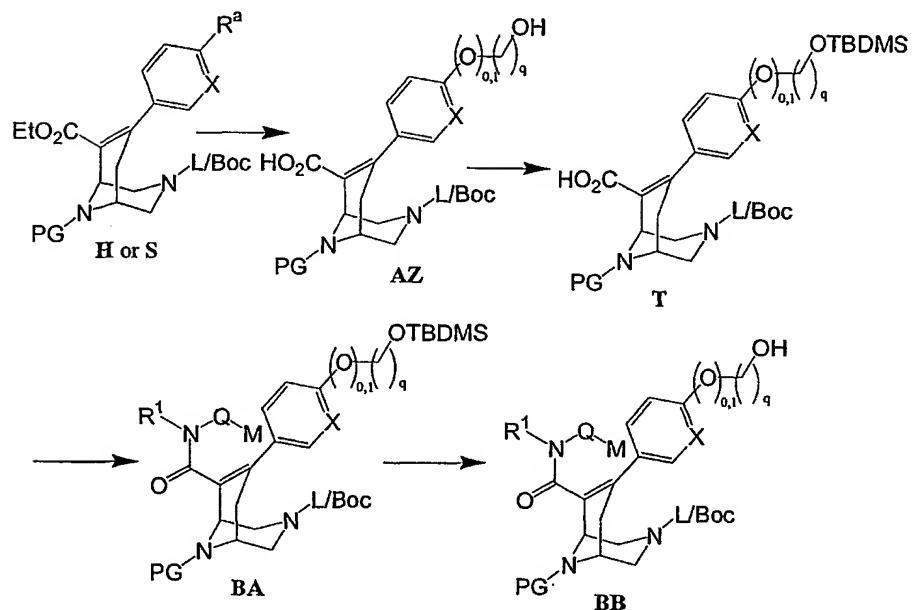
5 Alternatively, as shown in Scheme 15, AX might be transformed into a compound of type BL, than might be then lead to a compound of type AJ.

Scheme 15



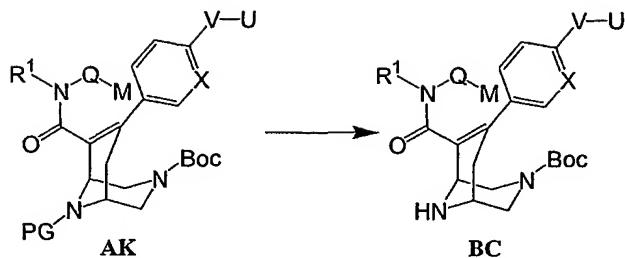
5 As shown in Scheme 16 compounds of type **H** or **S** might be transformed into compounds of type **AZ** (the substituent at the N(3) position be L or Boc). Then compounds of type **T** might be obtained that might be then transformed into compounds of type **BA**. Finally precursors of type **BB** might be prepared.

Scheme 16



Also, as shown in Scheme 17, compounds of type **AK** might be transformed into precursors of type **BC**.

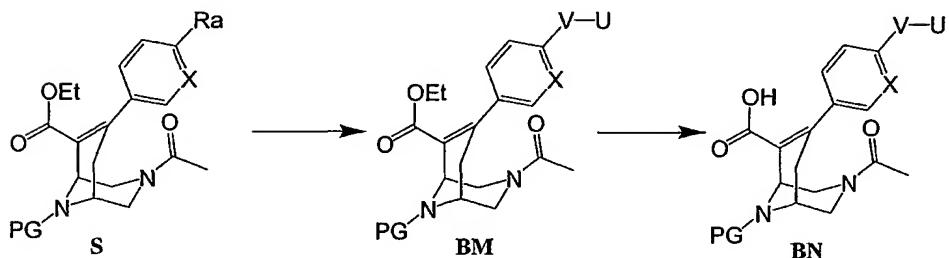
Scheme 17



- 5 Also, a compound of type S might be transformed into a compound of type BM, as shown in Scheme 18. A compound of type BM might be then saponified into a precursor of type BN.

Scheme 18

10

Preparation of the secondary amines

- 15 It might be necessary to prepare secondary amines as well. This might be done by reductive amination from the corresponding amine and aldehyde, or by amide coupling, from the corresponding amine and carboxylic acid, followed by reduction with LAH or borane. These standard procedures are well-described in the literature. (2-Allylphenyl)cyclopropylamine, necessary for instance in Scheme
20 13, might be prepared by allylation of 2-bromobenzaldehyde, protected as an acetal; subsequent deprotection to the 2-allylbenzaldehyde and reductive amination would lead to the desired amine.

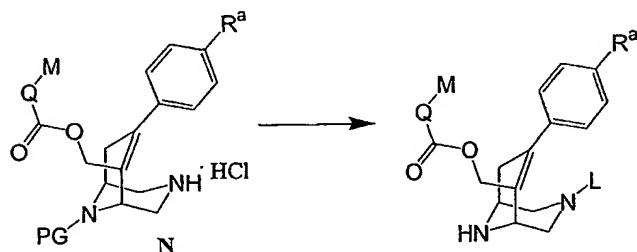
Preparation of final compounds

From precursors prepared as described above, the final compounds can be prepared using parallel chemistry techniques. For the specific examples, see the experimental part.

5

Diazabicyclononenes of type of N can be acylated, or alkylated, or sulfonated, using standard procedures (Scheme 19), and then directly deprotected to yield the final compounds (for numbering, see specific examples). In each case, purification by preparative HPLC might give the corresponding TFA salts or 10 formate salts.

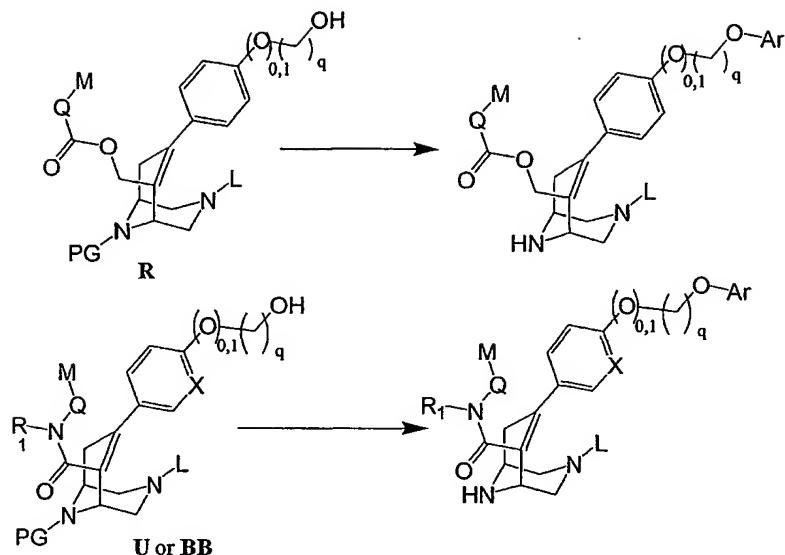
Scheme 19



15

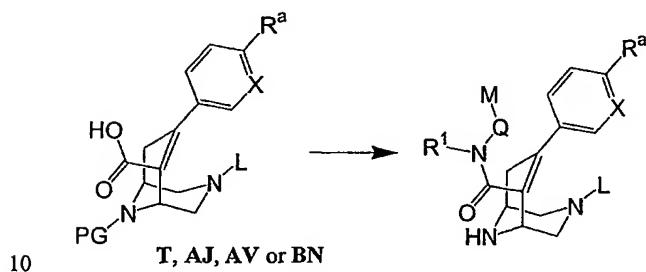
Precursors **R**, **U** or **BB** (with a L-substituent at the N(3)-position) might be transformed into the corresponding aryl ethers (Scheme 20), using the *Mitsunobu* reaction conditions. After deprotection, the final compounds are obtained.

Scheme 20



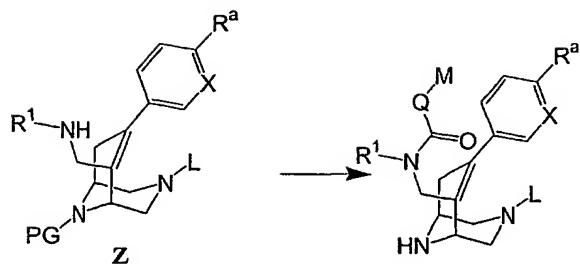
5 Precursors T, AJ or AV might be submitted to an amide coupling (Scheme 21). Deprotection would lead to the desired final compounds.

Scheme 21



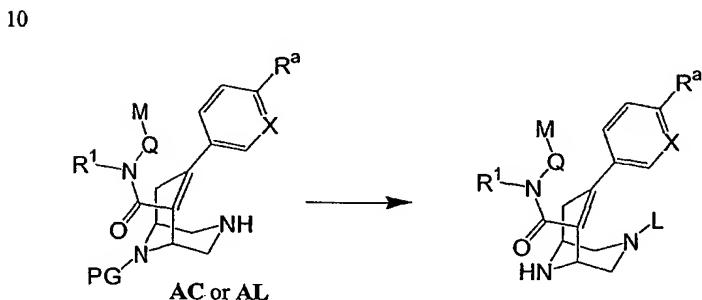
Compounds Z might be reacted with acylating (or sulfonating) reagents to lead to the corresponding amides (or sulfonamides) as well (Scheme 22). Deprotection would lead to the final compounds.

Scheme 22



5 Compounds of type **AC** or **AL** can be reacted as well with acylating, sulfonating or alkylating reagents (Scheme 23). After deprotection, the final compounds would be obtained.

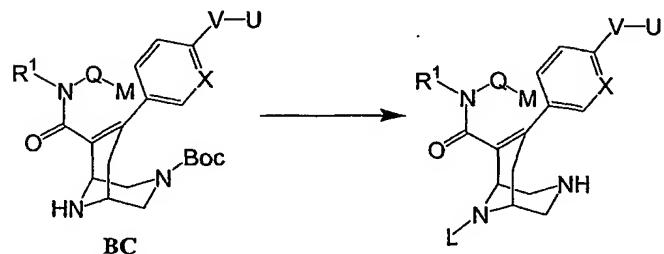
Scheme 23



Precursors of type **BC** might also lead to final compounds as indicated in Scheme 24.

15

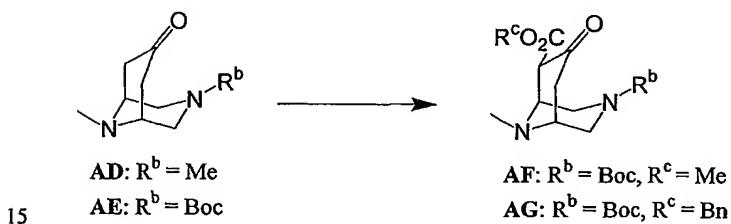
Scheme 24



Enantioselective synthesis:

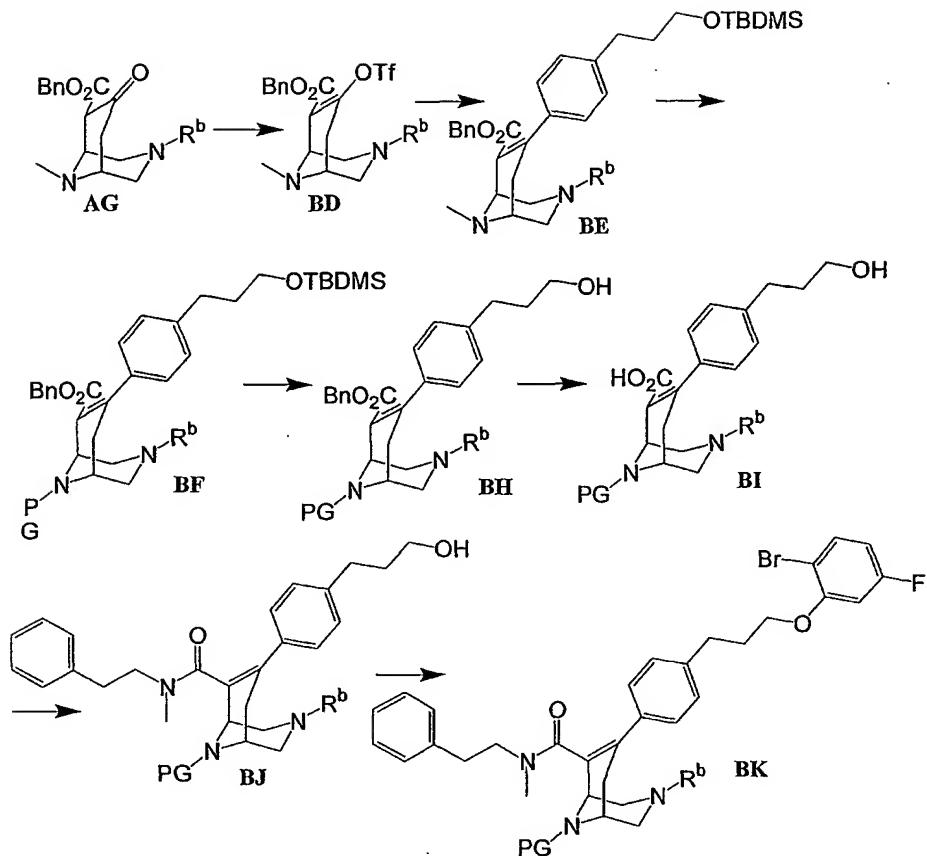
The compounds of the present invention contain at least two chiral centers which, however, are not independent from each other. The synthetic methods presented so far might lead to racemates. Both enantiomers might be prepared selectively starting from a *meso*-bicyclononane derivative, like compound AD (Blount B. K., Robinson, R., *J. Chem. Soc.*, 1932, 2485) or AE, prepared similarly to compound A with a subsequent decarboxylation (Scheme 25). For instance, compound AE might be stereoselectively acylated to bicyclononane derivatives AF or AG as already described elsewhere for similar compounds (Majewski M., Lasny R., *J. Org. Chem.*, 1995, 60, 5825). Similarly, the other enantiomer might be prepared.

Scheme 25



Finally precursor BK might be prepared as described in Scheme 26.

Scheme 26



5 The compounds of formula I and their pharmaceutically acceptable acid addition salts can be used as medicaments, e. g. in the form of pharmaceutical preparations for enteral, parenteral, or topical administration. They can be administered, for example, perorally, e. g. in the form of tablets, coated tablets, dragées, hard and soft gelatine capsules, solutions, emulsions or suspensions, rectally, e. g. in the
10 form of suppositories, parenterally, e. g. in the form of injection solutions or infusion solutions, or topically, e. g. in the form of ointments, creams or oils.

The production of pharmaceutical preparations can be effected in a manner which will be familiar to any person skilled in the art by bringing the described
15 compounds of formula I and their pharmaceutically acceptable acid addition salts,

optionally in combination with other therapeutically valuable substances, into a galenical administration form together with suitable, non-toxic, inert, therapeutically compatible solid or liquid carrier materials and, if desired, usual pharmaceutical adjuvants.

5

Suitable carrier materials are not only inorganic carrier materials, but also organic carrier materials. Thus, for example, lactose, corn starch or derivatives thereof, talc, stearic acid or its salts can be used as carrier materials for tablets, coated tablets, dragées and hard gelatine capsules. Suitable carrier materials for soft 10 gelatine capsules are, for example, vegetable oils, waxes, fats and semi-solid and liquid polyols (depending on the nature of the active ingredient no carriers are, however, required in the case of soft gelatine capsules). Suitable carrier materials for the production of solutions and syrups are, for example, water, polyols, sucrose, invert sugar and the like. Suitable carrier materials for injections are, for 15 example, water, alcohols, polyols, glycerols and vegetable oils. Suitable carrier materials for suppositories are, for example, natural or hardened oils, waxes, fats and semi-liquid or liquid polyols. Suitable carrier materials for topical preparations are glycerides, semi-synthetic and synthetic glycerides, hydrogenated oils, liquid waxes, liquid paraffins, liquid fatty alcohols, sterols, polyethylene 20 glycols and cellulose derivatives.

Usual stabilizers, preservatives, wetting and emulsifying agents, consistency-improving agents, flavour-improving agents, salts for varying the osmotic pressure, buffer substances, solubilizers, colorants and masking agents and 25 antioxidants come into consideration as pharmaceutical adjuvants.

The dosage of compounds of formula I can vary within wide limits depending on the disease to be controlled, the age and the individual condition of the patient and the mode of administration, and will, of course, be fitted to the individual 30 requirements in each particular case. For adult patients a daily dosage of about 1 mg to about 1000 mg, especially about 50 mg to about 500 mg, comes into consideration.

The pharmaceutical preparations conveniently contain about 1 - 500 mg, preferably 5 - 200 mg of a compound of formula I.

5 The following examples serve to illustrate the present invention in more detail. They are, however, not intended to limit its scope in any manner.

Examples

10 General remarks

The compounds were characterized at least by LC-MS and $^1\text{H-NMR}$. Only the LC-MS data are given here.

15 Abbreviations

AcCl	Acetyl chloride
ACE	Angiotensin Converting Enzyme
AcOH	Acetic acid
20 Ang	Angiotensin
aq.	aqueous
9-BBN	9-Borabicyclo[3.3.1]nonane
Bn	Benzyl
Boc	<i>tert</i> -Butyloxycarbonyl
25 BSA	Bovine serum albumine
BuLi	<i>n</i> -Butyllithium
CDI	1,1-Carbonyldiimidazol
conc.	concentrated
DIBAL	Diisobutylaluminium hydride
30 DIPEA	Diisopropylethylamine
DMAP	4- <i>N,N</i> -Dimethylaminopyridine
DMF	<i>N,N</i> -Dimethylformamide

	DMSO	Dimethylsulfoxide
	EDC·HCl	Ethyl- <i>N,N</i> -dimethylaminopropylcarbodiimide hydrochloride
	EIA	Enzyme immunoassay
	eq.	equivalent
5	Et	Ethyl
	EtOAc	Ethyl acetate
	FC	Flash Chromatography
	HOBt	Hydroxybenzotriazol
	KHMDS	Potassium hexamethyldisilazide
10	LAH	Lithium aluminium hydride
	MeOH	Methanol
	MPLC	Medium Pressure Liquid Chromatography
	NMO	<i>N</i> -Methylmorpholine <i>N</i> -oxide
	org.	organic
15	PG	protecting group
	Ph	Phenyl
	RAS	Renin Angiotensin System
	RP18	Reversed phase column, filled with C ₁₈ hydrocarbon
	rt	room temperature
20	SEM	Trimethylsilylethoxymethyl
	sol.	Solution
	TBAF	Tetra- <i>n</i> -butylammonium fluoride
	TBDMS	<i>tert</i> -Butyldimethylsilyl
	TBDPS	<i>tert</i> -Butyldiphenylsilyl
25	tBuOH	<i>tert</i> -Butanol
	tBuOK	Potassium <i>tert</i> -butylate
	Tf	Trifluoromethylsulfonyl
	TFA	Trifluoroacetic acid
	THF	Tetrahydrofuran
30	TLC	Thin Layer Chromatography
	TMAD	<i>N,N,N',N'</i> -Tetramethylazodicarboxamide

Preparation of the precursors**(*rac.*)-(1*R*^{*, 5S^{*})-9-Methyl-7-oxo-3,9-diazabicyclo[3.3.1]nonane-3,6-dicarboxylic acid 3-*tert*-butyl ester 6-ethyl ester (A)}**

5 (4-Benzyl-6-ethoxycarbonylmethyl-1-methyl-piperazin-2-yl)acetic acid ethyl ester (Patent WO 92/05174) (71.0 g, 0.196 mol) was dissolved in MeOH (400 mL). TFA (77.8 mL, 1.02 mol) was added and the flask was purged with nitrogen. Pd/C (10%, 50% moisture, 3.6 g) was added. The flask was closed and purged
10 with hydrogen (3x). After 1 day, the mixture was filtered through *Celite* and washed with MeOH. The solvents were removed under reduced pressure and the foamy residue (92.7 g) was dried under high vacuum. A sol. of tBuOK (117.2 g, 1.04 mol) in toluene (3.07 L) was heated to reflux. A sol. of the crude piperazine formerly obtained, dissolved in THF (300 mL), was added dropwise over 50 min.
15 The black mixture was stirred for 10 further min. and allowed to cool to rt. The mixture was cooled to 0 °C and AcOH (36.6 mL, 0.635 mol) was added. The solvents were removed under reduced pressure. This crude material was suspended in CH₂Cl₂ (400 mL) and cooled to 0 °C. DIPEA (19.1 mL, 112 mmol) was added. A sol. of Boc₂O (24.3 g, 113 mmol) in CH₂Cl₂ (200 mL) was added
20 dropwise. The mixture was stirred for 1 h at 0 °C, then 1 h at rt. The mixture was washed with aq. 10% Na₂CO₃ (2x). The org. extracts were dried over MgSO₄, filtered and the solvents were evaporated under reduced pressure. The residue was purified by FC (EtOAc/heptane 1:1 → EtOAc). The title compound was obtained as oil (24.5 g, 38%). R_f = 0.05 (EtOAc/heptane 1:1) or 0.56
25 (MeOH/CH₂Cl₂ 1:9). LC-MS: R_t = 2.94; ES+: 325.19.

(*rac.*)-(1*R*^{, 5S^{*})-9-Methyl-7-trifluoromethanesulfonyloxy-3,9-diazabicyclo-[3.3.1]non-6-ene-3,6-dicarboxylic acid 3-*tert*-butyl ester 6-ethyl ester (B)}

30 A sol. of bicyclononanone A (2.22 g, 6.80 mmol) in THF (50 mL) was cooled to 0 °C and NaH (about 60% in mineral oil, 326 mg, about 8.2 mmol) was added. A gas evolution was observed. After 20 min, Tf₂NPh (3.22 g, 9.00 mmol) was

added. 10 min later, the ice bath was removed. After 3 h, the sol. was diluted with EtOAc and washed with brine (1x). The org. extracts were dried over MgSO₄, filtered, and the solvents were removed under reduced pressure. Purification by FC (EtOAc/heptane 3:1 → EtOAc) yielded the title compound as
5 an oil (2.50 g, 80%). R_f = 0.15 (EtOAc/heptane 1:1). LC-MS: R_t = 4.73; ES+: 458.95.

Compounds of type C

10 3-(4-Bromophenyl)prop-1-yl 2-chlorophenyl ether (C1)

To a sol. of 3-(4-bromophenyl)propan-1-ol (Glover S. A., *et al.*; *Tetrahedron*, 1990, 46, 7247; 24.5 g, 0.114 mol) in toluene (600 mL) under nitrogen were added 2-chlorophenol (17.4 mL, 0.171 mmol), diisopropyl azodicarboxylate (33.1 mL, 0.171 mol) and tri-*n*-butylphosphine (42.2 mL, 0.171 mol). The sol. was heated to reflux and stirred under reflux overnight. The sol. was allowed to cool to rt and the solvents were removed under reduced pressure. The residue was diluted in EtOAc and washed with aq. 1M HCl (1x) and aq. 1M NaOH (2x). The org. extracts were dried over MgSO₄, filtered and the solvents were removed under reduced pressure. Purification of the residue by FC (petroleum ether → Et₂O/petroleum ether 1:99 → 1:19) led to the title compound (15.1 g, 41%). R_f = 0.70 (EtOAc/heptane 1:3).

25 2-(4-Bromophenyl)eth-1-yl 2,3,5-trimethylphenyl ether (C2)

A mixture of 2-(4-bromophenyl)ethanol (20.0 mL, 143 mmol), 2,3,5-trimethylphenol (31.1 g, 229 mmol), azodicarboxylic dipiperidine (72.1 g, 286 mmol) and tributylphosphine (88 mL; 357 mmol) in toluene (2.00 L) was heated to reflux for 2 h. The mixture was allowed to cool to rt. The mixture was filtered, 30 washed with toluene and the solvents were partially removed under reduced pressure. The residue was diluted with Et₂O and washed with aq. 1M NaOH (2x). The org. extracts were dried over MgSO₄, filtered, and the solvents were removed

under reduced pressure. Purification of the residue by FC (petroleum ether → Et₂O/petroleum ether 1:3) yielded the title compound (33.1 g, 73%). LC-MS: R_f = 6.95.

5 1-Bromo-4-[3-(2-methoxybenzyl)propoxy]benzene (E)

4-Bromophenol (4.32 g, 25.0 mmol) and 1-(3-chloropropoxymethyl)-2-methoxybenzene (Vieira E., et al., *Bioorg. Med. Chem. Letters*, 1999, 9, 1397, 4.88 g, 22.7 mmol) were dissolved in DMF (150 mL). NaI (1.50 g, 0.10 mmol) and Cs₂CO₃ (16.3 g, 50.0 mmol) were added. The mixture was heated to 80 °C and stirred for 6 h, before it was allowed to cool to rt. After dilution with EtOAc (600 mL) the mixture was washed with water (1x), aq. 1M NaOH (1x), aq. 1M HCl (1x). The org. extracts were dried over MgSO₄ and filtered. The solvents were removed under reduced pressure. Purification of the residue by FC (Et₂O/petroleum ether 1:9 → 1:4) yielded the title compound (5.66 g, 71%). R_f = 0.60 (Et₂O/heptane 1:1).

5-Bromo-2-[3-(2-methoxybenzyloxy)propoxy]pyridine (F)

20 3-(5-Bromopyridin-2-yloxy)propan-1-ol (Patent Application WO 98/39328, 1.05 g, 4.51 mmol) was diluted at rt in DMF (24 ml) and the sol. cooled to 0 °C. NaH (55 – 65 weight % in mineral oil, 193 mg, 4.42 – 5.23 mmol) was added and the yellow mixture was stirred for 20 min. 2-Methoxybenzyl chloride (1.49 ml, 10.7 mmol) was added and the solution was allowed to warm to rt and was stirred 25 for 4 h. The mixture was quenched with ice and diluted with EtOAc (20 ml), washed with brine and water, dried over MgSO₄ and filtered. The solvents were evaporated under reduced pressure. Purification of the residue by FC (Et₂O, heptane 1:39 → 1:19) yielded the title compound (627 mg, 40 %) as an oil. R_f = 0.07 (Et₂O / heptane, 1:3).

(*rac.*)-(1*R*^{*}, 5*S*^{*})-9-Methyl-7-[4-(2-trimethylsilanylethoxymethoxy)phenyl]-3,9-diazabicyclo[3.3.1]non-6-ene-3,6-dicarboxylic acid 3-*tert*-butyl ester 6-ethyl ester (G1)

- 5 A sol. of [2-(4-bromophenoxymethoxy)ethyl]trimethylsilane (Blass B. E., *et al.*, *Tetrahedron Lett.*, 2001, 42, 1611, 4.13 g, 13.6 mmol) in THF (30 mL) was cooled to -78 °C. BuLi (1.6 M in hexane, 9.1 mL, 14.6 mmol) was added. The sol. was stirred at -78 °C for 30 min. A sol. of ZnCl₂, prepared from ZnCl₂ (2.23 g, about 16.4 mmol) dried under high vacuum for 2 h at 140 °C and THF (35 mL),
- 10 was added and the resulting sol. was allowed to warm up to rt. A sol. of bicyclononene B (2.50 g, 5.45 mmol) in THF (5 mL) and then Pd(PPh₃)₄ (157 mg, 0.136 mmol) were added. After 10 min, the reaction mixture was heated to reflux. After 90 min, the reaction mixture was allowed to cool to rt and quenched with aq. 1M HCl. The mixture was diluted with EtOAc and washed with aq. 10% Na₂CO₃ (1x). The org. extracts were dried over MgSO₄ and filtered. The solvents were removed under reduced pressure. Purification of the residue by FC (MeOH/CH₂Cl₂ 1:39 → 1:24 → 1:20) yielded the title compound as an oil (2.90 g, 99%). R_f = 0.39 (MeOH/CH₂Cl₂ 1:9). LC-MS: R_t = 4.35; ES+: 533.29.
- 15
- 20 (*rac.*)-(1*R*^{*}, 5*S*^{*})-7-{4-[2-(*tert*-Butyldimethylsilanyloxy)ethyl]phenyl}-9-methyl-3,9-diazabicyclo[3.3.1]non-6-ene-3,6-dicarboxylic acid 3-*tert*-butyl ester 6-ethyl ester (G2)

A sol. of [2-(4-bromophenyl)ethoxy]-*tert*-butyldimethylsilane (Fuji K., *et al.*, *Tetrahedron Lett.*, 1990, 31, 6553, 21.8 g, 69.1 mmol) in THF (250 mL) was cooled to -78 °C. BuLi (1.55M in hexane, 44.6 mL, 69.1 mmol) was added. The sol. turned temporarily orange, then yellowish. After 30 min, ZnCl₂ (1M in THF, 70 mL, 70 mmol, prepared as described for G1) was added. The sol. was allowed to warm up to rt. Vinyl triflate B (12.91 g, 28.2 mmol) dissolved in THF (20 mL),

25 and Pd(PPh₃)₄ (600 mg, 0.519 mmol) were added. The sol. was stirred at rt for 90 min and aq. 1M HCl (1 mL) was then added. The mixture was diluted in EtOAc and washed with aq. 1M NaOH (1x). The org. extracts were dried over MgSO₄,

30

filtered, and the solvents were removed under reduced pressure. Purification of the residue by FC (MeOH/CH₂Cl₂ 1:49 → 1:24 → 3:47 → 2:25) yielded the title compound (10.91 g, 71%). R_f = 0.65 (MeOH/CH₂Cl₂ 1:9). LC-MS: R_t = 5.32; ES+: 545.49.

5

(rac.)-(1*R*^{*}, 5*S*^{*})-7-{4-[3-(*tert*-Butyldimethylsilyloxy)propyl]phenyl}-9-methyl-3,9-diazabicyclo[3.3.1]non-6-ene-3,6-dicarboxylic acid 3-*tert*-butyl ester 6-ethyl ester (G3)

10 A sol. of [3-(4-bromophenyl)propoxy]-*tert*-butyldimethylsilane (Kiesewetter D. O., *Tetrahedron Asymmetry*, 1993, 4, 2183, 22.60 g, 68.6 mmol) in THF (250 mL) was cooled to -78 °C. BuLi (1.55M in hexane, 44.3 mL, 68.6 mmol) was added. The sol. turned orange, then dark green. After 30 min, ZnCl₂ (1M in THF, 69 mL, 69 mmol, prepared as described for G2) was added, whereas the sol. turned deep 15 yellow. The mixture was allowed to warm up to rt. Vinyl triflate B (12.91 g, 28.2 mmol) in THF (20 mL) and then Pd(PPh₃)₄ (600 mg, 0.519 mmol) were added. The mixture was stirred at rt for 90 min and aq. 1M HCl (1 mL) was added. The mixture was diluted with EtOAc and washed with aq. 1M NaOH (1x). The org. extracts were dried over MgSO₄, filtered and the solvents were removed under 20 reduced pressure. Purification of the residue by FC (MeOH/CH₂Cl₂ 1:49 → 1:24 → 3:47 → 2:23) yielded the title product (10.76 g, 70%). R_f = 0.60 (MeOH/CH₂Cl₂ 1:9). LC-MS: R_t = 4.95; ES+: 559.51.

25 **(rac.)-(1*R*^{*}, 5*S*^{*})-7-{6-[3-(2-Methoxybenzyl)propoxy]pyridin-3-yl}-9-methyl-3,9-diazabicyclo[3.3.1]non-6-ene-3,6-dicarboxylic acid 3-*tert*-butyl ester 6-ethyl ester (G4)**

30 A sol. of bromopyridinyl derivative F (300 mg, 852 μmol) in THF (10 ml) was cooled to -78 °C. BuLi (1.55M in hexane, 0.580 ml, 889 μmol) was added and the mixture was stirred for 30 min. ZnCl₂ (1M in THF, 0.94 ml, 0.94 mmol, prepared as described for G2) was added and the reaction mixture was allowed to

warm up to rt. Vinyl triflate **B** (259 mg, 596 μ mol) in THF (1 mL), was added, followed by Pd(PPh₃)₄ (20.4 mg, 16.6 μ mol). The mixture was refluxed for 2 h. The reaction was terminated upon addition of ice. After dilution with EtOAc (125 mL), the reaction mixture was washed with 10% aq. Na₂CO₃ and the org. extracts were dried over MgSO₄ and filtered. The solvents were evaporated under reduced pressure. Purification of the residue by FC (CH₂Cl₂/MeOH: 39:1 \rightarrow 29:1 \rightarrow 24:1 \rightarrow 19:1 \rightarrow 9:1) led to title compound (197 mg, 54 %). R_f = 0.35 (CH₂Cl₂/MeOH 9:1). LC-MS: R_t = 4.06; ES+: 582.78.

10 (*rac.*)-(1*R*^{*}, 5*S*^{*})-7-{4-[3-(2-Methoxybenzyloxy)propoxy]phenyl}-9-methyl-
3,9-diazabicyclo[3.3.1]non-6-ene-3,6-dicarboxylic acid 3-*tert*-butyl ester 6-
ethyl ester (**G5**)

A sol. of bromophenyl derivative **E** (5.60 g, 15.9 mmol) in THF (50 mL) was
15 cooled to -78 °C. BuLi (1.55M in hexane, 10.3 mL, 15.9 mmol) was added. The
sol. was stirred at -78 °C for 30 min and ZnCl₂ (1M in THF, 17.5 mL, 17.5 mmol,
prepared as described for **G2**) was added. After warming up to rt, a sol. of vinyl
triflate **B** (3.63 g, 7.90 mmol) in THF (5 mL), followed by Pd(PPh₃)₄ (205 mg,
0.177 mmol), were added. The mixture was heated to reflux while turning black.
20 After 1 h, the reaction mixture was allowed to cool to rt. Ice was added and the
mixture was diluted in EtOAc. The org. extracts were washed with aq. 1M NaOH
(1x) and dried over MgSO₄. The solvents were removed under reduced pressure.
Purification of the residue by FC (MeOH/CH₂Cl₂ 1:49 \rightarrow 3:97 \rightarrow 1:24 \rightarrow 1:19 \rightarrow
1:9) yielded the title compound (4.57 g, 99%). R_f = 0.50 (MeOH/CH₂Cl₂ 1:9).
25 LC-MS: R_t = 4.17; ES+: 581.60.

(*rac.*)-(1*R*^{*}, 5*S*^{*})-7-{4-[3-(2-Chlorophenoxy)propyl]phenyl}-9-methyl-3,9-di-
azabicyclo[3.3.1]non-6-ene-3,6-dicarboxylic acid 3-*tert*-butyl ester 6-ethyl
ester (**G6**)

30

A sol. of bromophenyl derivative **C1** (16.0 g, 49.0 mmol) in THF (700 mL) was
cooled to -78 °C. BuLi (1.55M in hexane, 34.8 mL, 54.0 mmol) was added. The

sol. was stirred at -78 °C for 30 min and ZnCl₂ (1M in THF, 54.0 mL, 54.0 mmol, prepared as described for G2) was added. After warming up to rt, a sol. of vinyl triflate B (15.0 g, 32.7 mmol) in THF (50 mL), followed by Pd(PPh₃)₄ (945 mg, 0.818 mmol), were added. The sol. was heated to reflux. After 30 min, the reaction mixture was allowed to cool to rt. Ice was added and the mixture was diluted in EtOAc. The org. phase was washed with aq. 1M NaOH (1x) and dried over MgSO₄. The solvents were removed under reduced pressure. Purification of the residue by FC (MeOH/CH₂Cl₂ 1:49 → 3:97 → 1:24 → 1:19 → 1:9) yielded the title compound (10.5 g, 58%). LC-MS: R_t = 4.41; ES+: 555.13.

10

(*rac.*)-(1*R*^{*}, 5*S*^{*})-7-{4-[2-(*tert*-Butyldimethylsilanyloxy)ethoxy]phenyl}-9-methyl-3,9-diazabicyclo[3.3.1]non-6-ene-3,6-dicarboxylic acid 3-*tert*-butyl ester 6-ethyl ester (G7)

15 BuLi (1.6 M in hexane, 218 mL, 350 mmol) was added to a sol. of [2-(4-bromo-phenoxy)ethoxy]-*tert*-butyldimethylsilane (Morita, C.; et al.al.; *Heterocycles*, 2000, 52, 1163; 129 g, 342 mmol) in THF (1.0 L) at -78 °C. The mixture was stirred for 1 h at -78 °C, and ZnCl₂ (1M in THF, 400 mL, 400 mmol, prepared as described for G2) was added. The mixture was allowed to warm up to rt. 20 Bicyclononene B (78.4 g, 171 mmol) and Pd(PPh₃)₄ (4.94 g, 4.28 mmol) were added. The mixture was heated to reflux for 0.5 h, and was allowed to cool to rt. Aq. 1M HCl (2 mL) was added. The mixture was diluted with EtOAc (2 L) and washed with aq. 1M NaOH (750 mL). The aq. extracts were extracted back with EtOAc (1x). The combined org. extracts were dried over MgSO₄, filtered, and the solvents were removed under reduced pressure. Purification of the residue by FC (MeOH/CH₂Cl₂ 1:49 → 1:24 → 3:47 → 2:23) yielded the title compound (87.7 g, 91%). R_f = 0.60 (MeOH/CH₂Cl₂ 1:9). LC-MS: R_t = 4.74; ES+: 561.41.

25 (rac.)-(1*R*^{*}, 5*S*^{*})-7-{4-[2-(*tert*-Butyldiphenylsilanyloxy)ethyl]phenyl}-9-methyl-3,9-diazabicyclo[3.3.1]non-6-ene-3,6-dicarboxylic acid 3-*tert*-butyl ester 6-ethyl ester (G8)

BuLi (1.5 M in hexane, 13.4 mL, 20 mmol) was added to a sol. [2-(4-bromo-phenyl)ethoxy]-*tert*-butyldiphenylsilane (8.79 g, 20.0 mmol, prepared from 2-(4-bromophenyl)ethanol, TBDS-Cl and imidazol in DMF) in THF (40 mL) at -78 °C. The mixture was stirred for 30 min at -78 °C, and ZnCl₂ (1M in THF, 24 mL, 5 24 mmol, prepared as described for G2) was added. The mixture was allowed to warm up to rt. Bicyclononene B (3.67 g, 8.00 mmol) and Pd(PPh₃)₄ (231 mg, 0.20 mmol) were added. The mixture was heated to 40 °C for 40 min, and was allowed to cool to rt. Aq. 1M HCl (2 mL) was added. The mixture was diluted with EtOAc and washed with aq. 1M NaOH. The aq. extracts were extracted back with 10 EtOAc (1x). The combined org. extracts were dried over MgSO₄, filtered, and the solvents were removed under reduced pressure. Purification of the residue by FC (MeOH/CH₂Cl₂ 1:49 → 1:24 → 3:47 → 2:23) yielded the title compound (4.32 g, 81%). LC-MS: R_t = 1.06; ES+: 669.49.

15 Compounds of type H

(*rac.*)-(IR*, 5S*)-7-[4-(2-Trimethylsilanylethoxymethoxy)phenyl]-3,9-diazabi-cyclo[3.3.1]non-6-ene-3,6,9-tricarboxylic acid 3-*tert*-butyl ester 6-ethyl ester 9-(2,2,2-trichloro-1,1-dimethylethyl) ester (H1)

20 β,β,β—Trichloro-*tert*-butyl chloroformate (6.60 g, 27.5 mmol) was added to a sol. of bicyclononene G1 (2.93 g, 5.50 mmol) in 1,2-dichloroethane (60 mL). The sol. was heated to reflux. After 3 h, the reaction mixture was allowed to cool to rt, and the solvents were removed under reduced pressure. Purification of the residue by 25 FC (EtOAc/heptane 1:4 → 1:3 → 2:3 → 1:1) yielded the title compound as an oil (3.31 g, 83%). R_f = 0.52 (EtOAc/heptane 1:1). LC-MS: R_t = 7.40; ES+: 742.52.

(*rac.*)-(IR*, 5S*)-7-{4-[2-(*tert*-Butyldimethylsilanyloxy)ethyl]phenyl}-3,9-di-azabicyclo[3.3.1]non-6-ene-3,6,9-tricarboxylic acid 3-*tert*-butyl ester 6-ethyl 30 ester 9-(2,2,2-trichloro-1,1-dimethylethyl) ester (H2)

As for the preparation of compound **H1**, from bicyclononene **G2** (10.91 g, 20.0 mmol), β,β,β -trichloro-*tert*-butyl chloroformate (24.0 g, 100 mmol), and 1,2-dichloroethane (210 mL). Purification of the residue FC (EtOAc/heptane 1:9 → 1:4 → 2:3) yielded the title compound (13.75 g, 94%). $R_f = 0.64$ (EtOAc/heptane 5 1:1). LC-MS: $R_t = 7.66$; ES+: 755.37.

(*rac.*)-(1*R*^{*, 5*S*^{*})-7-{4-[3-(*tert*-Butyldimethylsilyloxy)propyl]phenyl}-3,9-di-azabicyclo[3.3.1]non-6-ene-3,6,9-tricarboxylic acid 3-*tert*-butyl ester 6-ethyl ester 9-(2,2,2-trichloro-1,1-dimethylethyl) ester (**H3**)}

10

As for the preparation of compound **H1**, from bicyclononene **G3** (10.96 g, 19.6 mmol), β,β,β -trichloro-*tert*-butyl chloroformate (23.5 g, 98.1 mmol), and 1,2-dichloroethane (210 mL). Purification of the residue by FC (EtOAc/heptane 1:9 → 1:4 → 2:3) yielded the title compound (13.50 g, 92%). $R_f = 0.58$ 15 (EtOAc/heptane 1:1). LC-MS: $R_t = 7.79$; ES+: 769.49.

(*rac.*)-(1*R*^{*, 5*S*^{*})-7-{6-[3-(2-Methoxybenzyloxy)propoxy]pyridin3-yl}-3,9-di-azabicyclo[3.3.1]non-6-ene-3,6,9-tricarboxylic acid 3-*tert*-butyl ester 6-ethyl ester 9-(2,2,2-trichloro-1,1-dimethylethyl) ester (**H4**)}

20

As for the preparation of compound **H1**, from bicyclononene **G4** (373 mg, 0.642 mmol), β,β,β -trichloro-*tert*-butyl chloroformate (770 mg, 3.11 mmol), and 1,2-dichloroethane (8 mL). Purification of the residue by FC (EtOAc/heptane 1:9 → 1:4 → 1:3) yielded the title compound (382 mg, 77%). $R_f = 0.47$ (EtOAc/heptane 25 1:1). LC-MS: $R_t = 7.14$; ES+: 770.50.

(*rac.*)-(1*R*^{*, 5*S*^{*})-7-{4-[3-(2-Methoxybenzyloxy)propoxy]phenyl}-3,9-diazabi-cyclo[3.3.1]non-6-ene-3,6,9-tricarboxylic acid 3-*tert*-butyl ester 6-ethyl ester 9-(2,2,2-trichloro-1,1-dimethylethyl) ester (**H5**)}

30

As for the preparation of compound **H1**, from bicyclononene **G5** (4.57 g, 7.87 mmol), β,β,β -trichloro-*tert*-butyl chloroformate (9.44 g, 39.4 mmol), and 1,2-

dichloroethane (100 mL). Purification of the residue by FC (EtOAc/heptane 1:9 → 1:4 → 1:1) yielded the title compound (5.35 g, 88%). $R_f = 0.46$ (EtOAc/heptane 1:1).

5 **(rac.)-(1*R*^{*}, 5*S*^{*})-7-{4-[3-(2-Chlorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-3,6,9-tricarboxylic acid 3-*tert*-butyl ester 6-ethyl ester 9-(2,2,2-trichloro-1,1-dimethylethyl) ester (H6)**

10 As for the preparation of compound H1, from bicyclononene G6 (10.51 g, 18.9 mmol), β,β,β-trichloro-*tert*-butyl chloroformate (22.7 g, 94.7 mmol), and 1,2-dichloroethane (350 mL). Purification of the residue by FC (EtOAc/heptane 1:8 → 1:4 → 1:1) yielded the title compound (12.5 g, 88%).

15 **(rac.)-(1*R*^{*}, 5*S*^{*})-7-{4-[2-(*tert*-Butyldimethylsilyloxy)ethoxy]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-3,6,9-tricarboxylic acid 3-*tert*-butyl ester 6-ethyl ester 9-(2,2,2-trichloro-1,1-dimethylethyl) ester (H7)**

20 As for the preparation of compound H1, from bicyclononene G7 (87.7 g, 156 mmol), β,β,β-trichloro-*tert*-butyl chloroformate (188 g, 784 mmol), and 1,2-dichloroethane (1.75 L). Purification of the residue by FC (EtOAc/heptane 1:19 → 1:3) yielded the title compound (111 g, 95%). $R_f = 0.75$ (EtOAc/heptane 1:1). LC-MS: $R_t = 7.84$.

25 **(rac.)-(1*R*^{*}, 5*S*^{*})-7-{4-[2-(*tert*-Butyldiphenylsilyloxy)ethyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-3,6,9-tricarboxylic acid 3-*tert*-butyl ester 6-ethyl ester 9-(2,2,2-trichloro-1,1-dimethylethyl) ester (H8)**

30 As for the preparation of compound H1, from bicyclononene G8 (4.32 g, 6.46 mmol), β,β,β-trichloro-*tert*-butyl chloroformate (7.75 g, 32.3 mmol), and 1,2-dichloroethane (100 mL). Purification of the residue by FC (EtOAc/heptane 1:19 → 1:3) yielded the title compound (4.97 g, 90%). $R_f = 0.75$ (EtOAc/heptane 1:1). LC-MS: $R_t = 1.35$.

Compounds of type J

(*rac.*)-(1*R*^{*,} 5*S*^{*})-6-Hydroxymethyl-7-[4-(2-trimethylsilanylethoxymethoxy)-5 phenyl]-3,9-diazabicyclo[3.3.1]non-6-ene-3, 9-dicarboxylic acid 3-*tert*-butyl ester 9-(2,2,2-trichloro-1,1-dimethylethyl) ester (**J1**)

A sol. of bicyclononene **H1** (3.31 g, 4.58 mmol) in CH₂Cl₂ (60 mL) was cooled to -78 °C. DIBAL (1M in toluene, 10.1 mL, 10.1 mmol) was added. The sol. was stirred for 30 min at -78 °C and was then allowed to warm slowly. DIBAL (5 mL) was added again after 1.5 h (-65 °C). Later, DIBAL was added successively in 5 mL-portions until TLC displayed no more starting material. Ice and water were then added at -50 °C. The cold bath was removed and the mixture warmed slowly to rt. More CH₂Cl₂ was added and the mixture was washed with aq. 1M HCl. The aq. phase was extracted with CH₂Cl₂ (1x) and the combined org. extracts were dried over MgSO₄. After filtration and evaporation of the solvents under reduced pressure, the residue was purified by FC (EtOAc/heptane 1:4 → 1:3 → 2:3) to yield the title compound as an oil (1.89 g, 60%). R_f = 0.50 (EtOAc/heptane 1:1). LC-MS: R_t = 7.08; ES+: 661.38, 702.83.

20

(*rac.*)-(1*R*^{*,} 5*S*^{*})-7-{4-2-(*tert*-Butyldimethylsilanyloxy)ethyl}phenyl}-6-hydroxymethyl-3,9-diazabicyclo[3.3.1]non-6-ene-3,9-dicarboxylic acid 3-*tert*-butyl ester 9-(2,2,2-trichloro-1,1-dimethylethyl) ester (**J2**)

To a sol. of bicyclononene **H2** (1.57 g, 2.32 mmol) in CH₂Cl₂ (40 mL) at -78 °C DIBAL (1M in toluene, 5.80 mL, 5.80 mmol) was added. The sol. was stirred at -78 °C for 1 h. Ice was added, and the mixture was allowed to warm up to rt. More CH₂Cl₂ was added and the org. extracts were washed with aq. 1M HCl (1x), dried over MgSO₄ and filtered. The solvents were removed under reduced pressure. Purification of the residue by FC (EtOAc/heptane 1:3) yielded the title compound (868 mg, 59%). LC-MS: R_t = 7.38; ES+: 715.48.

(rac.)-(1R, 5S*)-7-[4-[3-(*tert*-Butyldimethylsilyloxy)propyl]phenyl]-6-hydroxymethyl-3,9-diazabicyclo[3.3.1]non-6-ene-3,9-dicarboxylic acid 3-*tert*-butyl ester 9-(2,2,2-trichloro-1,1-dimethylethyl) ester (J3)*

5 A sol. of bicyclononene **H3** (2.39 g, 3.20 mmol) in CH₂Cl₂ (55 mL) was cooled to -78 °C. DIBAL (1M in toluene, 8.00 mL, 8.00 mmol) was added and the mixture was stirred at -78 °C for 1 h. Ice was added, and the mixture was allowed to warm up to rt. More CH₂Cl₂ was added and the org. extracts were washed with aq. 1M HCl (1x), dried over MgSO₄ and filtered. The solvents were removed under reduced pressure. Purification of the residue by FC (EtOAc/heptane 1:3) yielded the title compound (1.34 g, 59%). LC-MS: R_t = 7.59; ES+: 727.54.

10

15 *(rac.)-(1R*, 5S*)-6-Hydroxymethyl-7-[6-[3-(2-methoxybenzyloxy)propoxy]-pyridin-3-yl]-3,9-diazabicyclo[3.3.1]non-6-ene-3,9-dicarboxylic acid 3-*tert*-butyl ester 9-(2,2,2-trichloro-1,1-dimethylethyl) ester (J4)*

A sol. of bicyclononene **H4** (345 mg, 0.447 mmol) in CH₂Cl₂ (10 mL) was cooled to -78 °C. DIBAL (1M in toluene, 1.92 mL, 1.92 mmol) was added. The mixture was stirred at -78 °C for 1 h and again two portions of DIBAL (0.50 mL, 0.50 mmol) were added successively. After 2 h, ice was added. The mixture was allowed to warm up to rt and was diluted with more CH₂Cl₂. The org. extracts were washed with aq. 10% Na₂CO₃ (2x), dried over MgSO₄ and filtered. The solvents were removed under reduced pressure. Purification of the residue by FC (EtOAc/heptane 1:4 → 1:3 → 2:3 → 3:1) yielded the title compound (122 mg, 37%). R_f = 0.36 (EtOAc/heptane 1:1). LC-MS: R_t = 6.51; ES+: 728.49.

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(rac.)-(1R, 5S*)-6-Hydroxymethyl-7-(4-hydroxyphenyl)-3,9-diazabicyclo[3.3.1]non-6-ene-3,9-dicarboxylic acid 3-*tert*-butyl ester 9-(2,2,2-trichloro-1,1-dimethylethyl) ester (K)*

30

To a sol. of bicyclononene **J1** (1.89 g, 2.79 mmol) in THF/MeOH (1:1, 20 mL), a sol. of conc. H₂SO₄ (0.100 mL) in MeOH (10 mL) was added. The mixture was

stirred for 3 h at rt. The reaction mixture was diluted with EtOAc, washed with brine (1x) and aq. sat. NaHCO₃ (1x). The org. extracts were dried over MgSO₄ and filtered. The solvents were removed under reduced pressure. Crystallization of the residue from EtOAc/heptane led to the title compound (1.03 g, 67%). R_f = 5 0.14 (EtOAc/heptane 1:1). LC-MS: R_t = 5.17; ES-: 547.06.

(rac.)-(1*R*^{*}, 5*S*^{*})-6-Hydroxymethyl-7-{4-[3-(2-methoxybenzyloxy)propoxy]-phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-3,9-dicarboxylic acid 3-*tert*-butyl ester 9-(2,2,2-trichloro-1,1-dimethylethyl) ester (L)

10

To a sol. of bicyclononene K (1.03 g, 1.87 mmol) in DMF (20 mL) were added NaI (280 mg, 1.87 mmol), Cs₂CO₃ (609 mg, 1.87 mmol) and then 1-(3-chloropropoxymethyl)-2-methoxybenzene (Vieira E., et al., *Bioorg. Med. Chem. Letters*, 1999, 9, 1397, 400 mg, 1.87 mmol). The mixture was stirred at 100 °C. 15 After 1.5 h, another portion of Cs₂CO₃ (609 mg, 1.87 mmol) and 1-(3-chloropropoxymethyl)-2-methoxybenzene chloride (400 mg, 1.87 mmol) were added to complete the reaction. After 1.5 h later, the mixture was allowed to cool to rt and diluted with EtOAc. The org. extracts were washed with brine (1x). The org. extracts were dried over MgSO₄ and filtered. The solvents were removed 20 under reduced pressure. Purification of the residue by FC (EtOAc/heptane 1:4 → 1:3 → 2:3) yielded the title compound (1.00 g, 73%). R_f = 0.35 (EtOAc/heptane 1:1). LC-MS: R_t = 6.54.

(rac.)-(1*R*^{*}, 5*S*^{*})-7-{4-[3-(2-Methoxybenzyloxy)propoxy]phenyl}-6-[2-(2-methoxyphenyl)acetoxymethyl]-3,9-diazabicyclo[3.3.1]non-6-ene-3,9-dicarboxylic acid 3-*tert*-butyl ester 9-(2,2,2-trichloro-1,1-dimethylethyl) ester (M)

To a sol. of bicyclononene L (500 mg, 0.687 mmol) in CH₂Cl₂ (10 mL) were added (2-methoxyphenyl)acetic acid (206 mg, 1.37 mmol), DMAP (cat. amount), 30 DIPEA (0.230 mL, 1.34 mmol) and EDC·HCl (134 mg, 0.700 mmol). The sol. was stirred at rt for 90 min, when a second portion of DIPEA (0.100 mL, 0.584 mmol) was added. After 3 h, the reaction mixture was diluted in more CH₂Cl₂

and washed with water (1x). The org. extracts were dried over MgSO₄, filtered, and the solvents were removed under reduced pressure. Purification of the residue by FC (EtOAc/heptane 1:4 → 1:3 → 2:3) yielded the title compound (495 mg, 82%). R_f = 0.42 (EtOAc/heptane 1:1). LC-MS: R_t = 7.33; ES+: 897.33.

5

(*rac.*)-(IR*, 5S*)-7-{4-[3-(2-Methoxybenzyloxy)propoxy]phenyl}-6-[2-(2-methoxyphenyl)acetoxymethyl]-3,9-diazabicyclo[3.3.1]non-6-ene-9-carboxylic acid 2,2,2-trichloro-1,1-dimethylethyl ester hydrochloride salt (N)

10 A mixture of CH₂Cl₂ (3 mL) and HCl/dioxane (4M, 1 mL) was slowly added to bicyclononene M (495 mg, 0.585 mmol) in an ice bath. The resulting sol. was stirred at 0 °C. After 1 h, HCl/dioxane (4M, 0.5 mL) was added, and 1 h later the ice bath was removed. After 75 min, the solvents were removed under reduced pressure and the residue dried under high vacuum. The resulting foam was
15 estimated to contain about 80% of the title compound according to LC-MS and was used without further purification. LC-MS: R_t = 4.97; ES+: 774.97.

(*rac.*)-(IR*, 5S*)-6-[2-(2-Methoxyphenyl)acetoxymethyl]-7-{6-[3-(2-methoxybenzyloxy)propoxy]pyridin-3-yl}-3,9-diazabicyclo[3.3.1]non-6-ene-3,9-dicarboxylic acid 3-*tert*-butyl ester 9-(2,2,2-trichloro-1,1-dimethylethyl) ester (O)

To a sol. of bicyclononene J4 (122 mg, 0.167 mmol) in CH₂Cl₂ (5 mL) were added (2-methoxyphenyl)acetic acid (50 mg, 0.328 mmol), DIPEA (0.126 mL, 0.740 mmol), DMAP (cat. amount) and EDC-HCl (34 mg, 0.173 mmol). The mixture was stirred at rt for 3 h, then diluted in CH₂Cl₂, and washed with water (1x). The org. extracts were dried over MgSO₄ and filtered. The solvents were removed under reduced pressure. Purification of the residue by FC (EtOAc/heptane 1:4 → 1:3 → 2:3 → 1:1) yielded the title compound (108 mg, 74%). LC-MS: R_t = 7.34; ES+: 876.54.

30

(rac.)-(1R, 5S*)-7-{6-[3-(2-Methoxybenzyloxy)propoxy]pyridin-3-yl}-6-[2-(2-methoxyphenyl)acetoxyethyl]-3,9-diazabicyclo[3.3.1]non-6-ene-9-carboxylic acid 2,2,2-trichloro-1,1-dimethylethyl ester dihydrochloride salt (P)*

5 Bicyclononene **O** (114 mg, 0.130 mmol) was dissolved in CH₂Cl₂ (2 mL). The sol. was cooled to -40 °C and 4M HCl/dioxane (2 mL) was added. The sol. was stirred for 50 min while warming up slowly to 0 °C and then stirred for 1.5 h at 0 °C. The solvents were rapidly removed under reduced pressure. The residue was dried under high vacuum to give the title compound (156 mg) as a foam that was
10 used without further purification. LC-MS: R_t = 4.86; ES+: 776.48.

Compounds of type Q

15 *(rac.)-(1R*, 5S*)-7-{4-[2-(tert-Butyldimethylsilyloxy)ethyl]phenyl}-6-[2-(2-methoxyphenyl)acetoxyethyl]-3,9-diazabicyclo[3.3.1]non-6-ene-3,9-dicarboxylic acid 3-tert-butyl ester 9-(2,2,2-trichloro-1,1-dimethylethyl) ester (Q1)*

To a sol. of bicyclononene **J2** (868 mg, 1.25 mmol) in CH₂Cl₂ (20 mL) were added (2-methoxyphenyl)acetic acid (343 mg, 2.06 mmol), DIPEA (0.652 mL, 3.81 mmol), DMAP (cat. amount), and EDC·HCl (201 mg, 1.05 mmol). The mixture was stirred at rt for 1 h and EDC·HCl (73 mg, 0.38 mmol) and DIPEA (0.163 mmol, 0.952 mmol) were added again. After 30 min. the reaction mixture was diluted with CH₂Cl₂, washed with aq. 1M HCl (1x), and aq. sat. NaHCO₃ (1x). The org. extracts were dried over MgSO₄ and filtered. The solvents were
20 removed under reduced pressure. Purification of the residue by FC (EtOAc/heptane 1:3) yielded the title compound (948 mg, 90%). LC-MS: R_t = 7.98; ES+: 861.51.

30 *(rac.)-(1R*, 5S*)-7-{4-[3-(tert-Butyldimethylsilyloxy)propyl]phenyl}-6-[2-(2-methoxyphenyl)acetoxyethyl]-3,9-diazabicyclo[3.3.1]non-6-ene-3,9-dicarboxylic acid 3-tert-butyl ester 9-(2,2,2-trichloro-1,1-dimethylethyl) ester (Q2)*

To a sol. of bicyclononene **J3** (1.34 g, 1.90 mmol) in CH₂Cl₂ (30 mL), were added 2-methoxyphenylacetic acid (633 mg, 3.81 mmol), DIPEA (0.652 mL, 3.81 mmol), DMAP (cat. amount), and EDC·HCl (402 mg, 2.09 mmol). The sol. was stirred at rt for 1 h and EDC·HCl (73 mg, 0.38 mmol) and DIPEA (0.163 mmol, 5 0.952 mmol) were added again. After 30 min., the reaction mixture was diluted with CH₂Cl₂, washed with aq. 1M HCl (1x) and aq. sat. NaHCO₃ (1x). The org. extracts were dried over MgSO₄ and filtered. The solvents were removed under reduced pressure. Purification of the residue by FC (EtOAc/heptane 1:3) yielded the title compound (1.47 g, 90%). LC-MS: R_t = 8.17; ES+: 875.53.

10

Compounds of type R

(*rac.*)-(1*R*^{*}, 5*S*^{*})-3-Acetyl-7-[4-(2-hydroxyethyl)phenyl]-6-[2-(2-methoxyphenyl)acetoxyethyl]-3,9-diazabicyclo[3.3.1]non-6-ene-9-carboxylic acid
15 2,2,2-trichloro-1,1-dimethylethyl ester (**R1**)

Bicyclononene **Q1** (948 mg, 1.13 mmol) was dissolved in CH₂Cl₂ (20 mL). The sol. was cooled to 0 °C and 4M HCl/dioxane (20 mL) was added. After 2.25 h the solvents were rapidly removed under reduced pressure, and the residue was 20 immediately dried under high vacuum. The resulting foam was then dissolved in THF (55 mL) and cooled to -78 °C. DIPEA (0.774 mL, 4.51 mmol) and DMAP (cat. amount) were added, followed by the addition of acetyl chloride (0.064 mL, 0.91 mmol). After 2.5 h at -78 °C, MeOH (20 mL) was added, and the reaction mixture was allowed to warm to rt. The reaction mixture was diluted in EtOAc, 25 washed with aq. 1 M HCl (1x), and the org. extract were dried over MgSO₄ and filtered. The solvents were removed under reduced pressure. Purification of the residue by FC (EtOAc/heptane 1:2) yielded the title compound (651 mg, 55%). LC-MS: R_t = 5.47; ES+: 689.05.

30 (*rac.*)-(1*R*^{*}, 5*S*^{*})-3-Acetyl-7-[4-(3-hydroxypropyl)phenyl]-6-[2-(2-methoxyphenyl)acetoxyethyl]-3,9-diazabicyclo[3.3.1]non-6-ene-9-carboxylic acid
2,2,2-trichloro-1,1-dimethylethyl ester (**R2**)

To a sol. of bicyclononene **Q2** (1.47 g, 1.72 mmol) in CH₂Cl₂ (25 mL), 4M HCl/dioxane (25 mL) was added at 0 °C. After 2.25 h at 0 °C the solvents were rapidly removed under reduced pressure, and the residue was immediately dried
5 under high vacuum. The resulting foam was then dissolved in THF (75 mL) and cooled to -78 °C. DIPEA (1.18 mL, 6.88 mmol) and DMAP (cat. amount) were added, followed by slow addition of acetyl chloride (0.098 mL, 1.38 mmol). After 2.5 h at -78 °C MeOH (80 mL) was added, and the reaction mixture was allowed to warm up to rt. After dilution with EtOAc, the mixture was washed with aq. 1
10 M HCl (1x) and the org. extracts were dried over MgSO₄ and filtered. The solvents were removed under reduced pressure. Purification of the residue by FC (EtOAc/heptane 1:2) yielded the title compound (651 mg, 55%).

Compounds of type S

15

(*rac*.)-(1*R*^{*}, 5*S*^{*})-3-Acetyl-7-{4-[3-(2-methoxybenzyloxy)propoxy]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6,9-dicarboxylic acid 6-ethyl ester 9-(2,2,2-trichloro-1,1-dimethylethyl) ester (**S1**)

20 A sol. of bicyclononene **H5** (5.35 g, 6.95 mmol) in CH₂Cl₂ (30 mL) was cooled to 0°C. 4M HCl/dioxane (30 mL) was added. The sol. was stirred at 0 °C for 3.5 h, the solvents were removed under reduced pressure and the residue dried at high vacuum. The resulting foam was dissolved in THF (100 mL) and cooled to -78 °C. DIPEA (5.80 mL, 34.7 mmol) was added. A sol. of acetyl chloride (0.494
25 mL, 6.95 mmol) in THF (10 mL) was added slowly. The reacton mixture was stirred at -78 °C for 90 min, then allowed allowed to warm to rt and diluted in MeOH (5 mL), then in EtOAc. The org. extracts were washed with aq. 1M HCl (2x), aq. sat. NaHCO₃ (1x), dried over MgSO₄, filtered and the solvents were removed under reduced pressure. Purification of the residue by FC (EtOAc/heptane 1:4 → 1:3 → 1:1 → EtOAc) yielded the title compound (3.67 g,
30 74%). R_f = 0.50 (EtOAc). LC-MS: 6.22; ES+: 711.31.

(rac.)-(1R, 5S*)-3-Acetyl-7-[4-[3-(2-chlorophenoxy)propyl]phenyl]-3,9-diazabicyclo[3.3.1]non-6-ene-6,9-dicarboxylic acid 6-ethyl ester 9-(2,2,2-trichloro-1,1-dimethylethyl) ester (S2)*

5 As for compound S1, from H6 (10.8 g, 14.5 mmol), CH₂Cl₂ (110 mL), 4M HCl/dioxane (110 mL), THF (220 mL), DIPEA (12.4 mL, 72.6 mmol), DMAP (89 mg, 0.73 mmol), acetyl chloride (1.24 mL, 17.4 mmol), and MeOH (5 mL). Purification of the residue by FC (EtOAc/heptane 1:3 → 1:1 → EtOAc) yielded the title compound (8.59 g, 86%). R_f = 0.43 (EtOAc). LC-MS: 6.32; ES+:
10 684.99.

(rac.)-(1R, 5S*)-3-Acetyl-7-[4-(2-hydroxyethyl)phenyl]-3,9-diazabicyclo[3.3.1]non-6-ene-6,9-dicarboxylic acid 6-ethyl ester 9-(2,2,2-trichloro-1,1-dimethylethyl) ester (S3)*

15 As for compound S1, from H2 (3.00 g, 4.08 mmol), CH₂Cl₂ (30 mL), 4M HCl/dioxane (30 mL), THF (60 mL), DMAP (25 mg, 0.204 mmol), DIPEA (2.74 mL, 16.4 mmol), acetyl chloride (0.343 mL, 4.08 mmol), and MeOH (5 mL). Purification of the residue by FC (EtOAc/heptane 1:1 → EtOAc → MeOH/EtOAc
20 1:9) led to the title compound (1.80 g, 79%). R_f = 0.20 (EtOAc).

(rac.)-(1R, 5S*)-3-Acetyl-7-[4-(2-hydroxyethoxy)phenyl]-3,9-diazabicyclo[3.3.1]non-6-ene-6,9-dicarboxylic acid 6-ethyl ester 9-(2,2,2-trichloro-1,1-dimethylethyl) ester (S4)*

25 As for compound S1, from H7 (5.80 g, 7.73 mmol), CH₂Cl₂ (60 mL), 4M HCl/dioxane (60 mL), THF (50 mL), DMAP (47 mg, 0.384 mmol), DIPEA (5.29 mL, 31.7 mmol), acetyl chloride (0.604 mL, 8.08 mmol), and MeOH (5 mL). Purification of the residue by FC (EtOAc/heptane 1:1 → EtOAc → MeOH/EtOAc
30 1:9) led to the title compound (3.07 g, 69%). R_f = 0.20 (EtOAc). LC-MS: R_t = 4.80; ES+: 576.93.

(rac.)-(1R, 5S*)-3-Acetyl-7-[4-(3-hydroxypropyl)phenyl]-3,9-diazabicyclo[3.3.1]non-6-ene-6,9-dicarboxylic acid 6-ethyl ester 9-(2,2,2-trichloro-1,1-dimethylethyl) ester (S5)*

5 As for compound **S1**, from **H3** (5.04 g, 6.74 mmol), CH₂Cl₂ (80 mL), 4M HCl/dioxane (80 mL), THF (80 mL), without DMAP, DIPEA (4.62 mL, 27.0 mmol), acetyl chloride (0.430 mL, 6.06 mmol), and MeOH (5 mL). Purification of the residue by FC (EtOAc/heptane 1:1 → EtOAc → MeOH/EtOAc 1:9) led to the title compound (3.23 g, 83%). R_f = 0.20 (EtOAc). LC-MS: R_t = 1.00; ES+:

10 575.13.

Compounds of type T

15 *(rac.)-(1R*, 5S*)-3-Acetyl-7-{4-[3-(2-methoxybenzyloxy)propoxy]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6,9-dicarboxylic acid 9-(2,2,2-trichloro-1,1-dimethylethyl) ester (T1)*

To a sol. of bicyclononene **S1** (3.67 g, 5.15 mmol) in EtOH (27 mL) was added aq. NaOH (1M, 27 mL, 27 mmol). The mixture was heated to 80 °C for 3 h and 20 then allowed to cool to rt. After adjustment of the pH to 1-2 with aq. 1M HCl and extraction with EtOAc (2x), the combined org. extracts were dried over MgSO₄, filtered, and evaporated under reduced pressure. Purification of the residue by FC (EtOAc/heptane 1:1 → 3:1 → EtOAc → MeOH/EtOAc 1:9) yielded the title compound (1.45 g, 41%). LC-MS: R_t = 5.50; ES+: 683.24.

25

(rac.)-(1R, 5S*)-3-Acetyl-7-{4-[3-(2-chlorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6,9-dicarboxylic acid 9-(2,2,2-trichloro-1,1-dimethylethyl) ester (T2)*

30 From bicyclononene **S2** (8.59 g, 12.5 mmol) the title compound (4.29 g, 52%) was obtained after purification by FC (MeOH/CH₂Cl₂ 1:99 → 1:49 → 3:97 → 1:24) as described for **T1**. LC-MS: R_t = 5.61; ES-: 655.24.

(*rac.*)-(1*R*^{*}, 5*S*^{*})-3-Acetyl-7-[4-(2-hydroxyethoxy)phenyl]-3,9-diazabicyclo[3.3.1]non-6-ene-6,9-dicarboxylic acid 9-(2,2,2-trichloro-1,1-dimethylethyl) ester (**T3**)

5

From bicyclononene **S4** (3.07 g, 5.73 mmol) in EtOH (119 mL) and aq. 1M NaOH (119 mL) the title compound (1.88 g, 60%) was obtained after purification by FC (MeOH/CH₂Cl₂ 1:99 → 1:49 → 3:97 → 1:24) as described for **T1**. LC-MS: R_t = 4.32; ES+: 548.96.

10

(*rac.*)-(1*R*^{*}, 5*S*^{*})-3-Acetyl-7-{4-[3-(*tert*-butyldimethylsilyloxy)propyl]-phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6,9-dicarboxylic acid 9-(2,2,2-trichloro-1,1-dimethylethyl) ester (**T4**)

15 A mixture of bicyclononene **AZ1** (1.60 g, 2.93 mmol), imidazol (797 mg, 11.7 mmol) and TBDMS-Cl (1.1 g, 7.30 mmol) in DMF (20 mL) was stirred at rt overnight. Aq. sat. NH₄Cl was added and the mixture was extracted with hexane (3x). The combined org. extracts were dried over MgSO₄, filtered, and the solvents were evaporated under reduced pressure. A mixture of this crude product 20 and K₂CO₃ (0.2 g) in THF (30 ml), MeOH (10 ml), and H₂O (10 ml) was stirred at rt for 3 h. Aq. sat. NH₄Cl was added and this mixture was extracted with Et₂O (3x). The combined org. extracts were dried over MgSO₄, filtered, and the solvents were removed under reduced pressure. This yielded the title compound 25 (1.85 g, 95%) that was used without further purification.

25

(*rac.*)-(1*R*^{*}, 5*S*^{*})-7-{4-[2-(*tert*-Butyldiphenylsilyloxy)ethyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-3,6,9-tricarboxylic acid 3-*tert*-butyl ester 9-(2,2,2-trichloro-1,1-dimethylethyl) ester (**T5**)

30 As for compound **T4**, but from bicyclononene **AZ2** (crude, about 5.79 mmol), imidazol (1.2 g, 17.6 mmol) and TBDPS-Cl (4.84 g, 17.6 mmol) in DMF (50 mL), then K₂CO₃ (0.5 g), THF (60 mL), MeOH (20 mL), and H₂O (20 mL). The crude

title compound (9.6 g, quantitative yield) was used further without purification.
LC-MS: $R_f = 1.26$.

Compounds of type U

5

(rac.)-(1R, 5S*)-3-Acetyl-7-(4-hydroxyphenyl)-6-(methylphenethylcarbamoyl)-3,9-diazabicyclo[3.3.1]non-6-ene-9-carboxylic acid 2,2,2-trichloro-1,1-dimethylethyl ester (U1)*

10 To a sol. of bicyclononene W1 (0.93 g, 1.37 mmol) in CH_2Cl_2 (10 mL) was added HCl/dioxane (10 mL) at 0 °C. After 15 min, the ice bath was removed. The reaction mixture was stirred at rt for 1 h and the solvents were removed under reduced pressure. After drying at high vacuum for 30 min., the resulting solid or foam was dissolved in THF (10 mL). DIPEA (0.983 mL, 5.48 mmol) and DMAP (cat. amount) were added. The sol. was cooled to -78 °C and a sol. of AcCl (0.0973 mL, 1.37 mmol) in THF (5 mL) was slowly added over 2 min. After 75 min at -78 °C MeOH (10 mL) was added and the mixture was allowed to warm up. After dilution in EtOAc, the reaction mixture was washed with aq. 1M HCl (1x) and aq. sat. NaHCO_3 (1x). The org. extracts were dried over MgSO_4 , filtered, 15 and evaporated under reduced pressure. Purification of the residue by FC (EtOAc/heptane 1:3 → 1:1 → 3:1 → EtOAc → MeOH/EtOAc 1:19 → 1:9) yielded the title compound (253 mg, 30%). $R_f = 0.30$ (EtOAc). LC-MS: $R_f = 5.12$; ES+: 622.31.

20
25 *(rac.)-(1R*, 5S*)-3-Acetyl-7-[4-(2-hydroxyethyl)phenyl]-6-(methylphenethylcarbamoyl)-3,9-diazabicyclo[3.3.1]non-6-ene-9-carboxylic acid 2,2,2-trichloro-1,1-dimethylethyl ester (U2)*

To a sol. of bicyclononene W2 (3.46 g, 4.88 mmol) in CH_2Cl_2 (35 mL) was added 30 HCl (4M in dioxane, 35 mL) at 0 °C. After 1 h the ice bath was removed, and stirring continued for 1 h at rt. The solvents were removed under reduced pressure and the residue dried under high vacuum. The resulting foam was

dissolved in THF (50 mL). DIPEA (3.34 mL, 19.5 mmol) and DMAP (cat. amount) were added. The reaction mixture was cooled to -78 °C and AcCl (0.347 mL, 4.88 mmol) in THF (20 mL) was added dropwise. After 2 h at -78 °C, AcCl (0.100 mL, 1.41 mmol) was added again, followed by a third portion of AcCl (0.050 mL, 0.71 mmol) 1.5 h later. MeOH (10 mL) was added after 30 min. and the mixture was allowed to warm up to rt. After dilution with EtOAc and washing with aq. 1M HCl (1x) and aq. sat. NaHCO₃ (1x), the org. extracts were dried over MgSO₄ and filtered. The solvents were removed under reduced pressure. Purification of the residue by FC (EtOAc/heptane 1:1 → EtOAc → MeOH/EtOAc 10 1:9) yielded the title compound as a colorless foam (2.06 g, 65%). R_f = 0.15 (EtOAc). LC-MS: R_t = 5.14; ES+: 650.21.

(rac.)-(1R, 5S*)-3-Acetyl-7-[4-(3-hydroxypropyl)phenyl]-6-(methylphenethylcarbamoyl)-3,9-diazabicyclo[3.3.1]non-6-ene-9-carboxylic acid 2,2,2-trichloro-1,1-dimethylethyl ester (U3)*

To a sol. of bicyclononene W3 (3.18 g, 4.40 mmol) in CH₂Cl₂ (30 mL) was added HCl (4M in dioxane, 30 mL) at 0 °C. After 1 h at 0 °C and 1 h at rt, the solvents were removed under reduced pressure and the residue dried under high vacuum. 20 The residue was dissolved in THF (45 mL), and DIPEA (3.02 mL, 17.6 mmol) and DMAP (cat. amount) were added. The sol. was cooled to -78 °C and a sol. of AcCl (0.313 mL, 4.40 mmol) in THF (15 mL) was added dropwise over 5 min. After 1.25 h, AcCl (0.070 mL, 0.984 mmol) was added again. After 30 min. MeOH (10 mL) was added and the mixture was allowed to warm up to rt. After 25 dilution in EtOAc and washing with aq. 1M HCl (1x) and aq. sat. NaHCO₃ (1x), the org. extracts were dried over MgSO₄ and filtered. The solvents were removed under reduced pressure. Purification of the residue by FC (EtOAc/heptane 1:1 → EtOAc → MeOH/EtOAc 1:9) yielded the title compound (2.92 g, 66%) as a foam. R_f = 0.23 (EtOAc). LC-MS: R_t = 5.24; ES+: 664.29.

**(*rac.*)-(IR*, 5S*)-3-Acetyl-7-[4-(2-hydroxyethoxy)phenyl]-6-(methylphenethylcarbamoyl)-3,9-diazabicyclo[3.3.1]non-6-ene-9-carboxylic acid
2,2,2-trichloro-1,1-dimethylethyl ester (U4)**

5 A mixture of bicyclononene T3 (1.88 g, 3.42 mmol), methylphenethylamine (1.49 mL, 10.3 mmol), DMAP (41 mg, 0.34 mmol), DIPEA (2.33 mL, 18.0 mmol), HOBt (46 mg, 0.34 mmol) and EDC·HCl (1.64 g, 8.55 mmol) in CHCl₃ (40 mL) was stirred overnight at rt. The mixture was diluted in CH₂Cl₂ and washed with aq. 1M HCl (2x) and aq. sat. NaHCO₃. The organic extracts were dried over 10 MgSO₄, filtered, and the solvents were removed under reduced pressure. Purification of the residue by FC (EtOAc/ heptane 1:4 → 1:1 → 4:1 → EtOAc) yielded the title compound (1.33 g, 58%). LC MS: R_t = 5.25; ES+: 666.08.

Compounds of type V

15

(*rac.*)-(IR*, 5S*)-7-(4-Hydroxyphenyl)-3,9-diazabicyclo[3.3.1]non-6-ene-3,6,9-tricarboxylic acid 3-*tert*-butyl ester 9-(2,2,2-trichloro-1,1-dimethylethyl) ester (V1)

20 To a sol. of bicyclononene H1 (4.18 g, 5.79 mmol) in EtOH (55 mL) aq. NaOH (1M, 55 mL, 55 mmol) was added. The mixture was stirred at 80 °C for 28 h before it was allowed to cool to rt and acidified to pH 1 with aq. 1M HCl. After extraction with EtOAc (3x) the combined org. extracts were dried over MgSO₄ and filtered. The solvents were removed under reduced pressure. Purification of 25 the residue by FC (MeOH/CH₂Cl₂ 1:49 → 3:97 → 1:24 → 1:19 → 1:9 → 1:4) yielded the title compound (1.50 g, 40%). R_f = 0.29. LC-MS: R_t = 4.91; ES-: 561.12.

30 (*rac.*)-(IR*, 5S*)-7-[4-(2-Hydroxyethyl)phenyl]-3,9-diazabicyclo[3.3.1]non-6-ene-3,6,9-tricarboxylic acid 3-*tert*-butyl ester 9-(2,2,2-trichloro-1,1-dimethyl-ethyl) ester (V2)

To a sol. of bicyclononene **H2** (13.75 g, 18.7 mmol) in EtOH (180 mL) aq. NaOH (1M, 180 mL, 180 mmol) was added. The mixture was stirred at 80 °C for 8 h and then left at -5 °C overnight. The mixture was acidified to pH 1 with aq. 1M HCl and extracted with EtOAc (3x). The org. extracts were dried over MgSO₄ 5 and filtered. The solvents were removed under reduced pressure. Purification of the residue by FC (MeOH/CH₂Cl₂ 1:49 → 3:47 → 1:24 → 1:19 → 1:9 → 1:4) yielded the title compound, contaminated with (*rac.*)-(1*R*^{*}, 5*S*^{*})-7-[4-(2-hydroxyethyl)phenyl]-3,9-diazabicyclo[3.3.1]non-7-ene-3,6,9-tricarboxylic acid 3-*tert*-butyl ester 9-(2,2,2-trichloro-1,1-dimethylethyl) ester (7.09 g, 64%). R_f 10 = 0.40 MeOH/CH₂Cl₂ 1:9). R_t = 4.90; ES-: 589.16.

(*rac.*)-(1*R*^{*}, 5*S*^{*})-7-[4-(3-Hydroxypropyl)phenyl]-3,9-diazabicyclo[3.3.1]non-6-ene-3,6,9-tricarboxylic acid 3-*tert*-butyl ester 9-(2,2,2-trichloro-1,1-dimethylethyl) ester (**V3**)

15 To a sol. of bicyclononene **H3** (13.50 g, 18.0 mmol) in EtOH (180 mL) aq. NaOH (1M, 180 mL, 180 mmol) was added. The mixture was heated to 40 °C and after 1 h to 80 °C. After 7 h, the mixture was left overnight at -5 °C. EtOH (100 mL) and aq. NaOH (1M, 50 mL, 50 mmol) were added and the sol. was heated to 80 20 °C for 6 h. After cooling to rt and adjustment of the pH to 1 with aq. 1M HCl, the mixture was extracted with EtOAc (3x). The combined org. extracts were dried over MgSO₄ and filtered. The solvents were removed under reduced pressure. Purification of the residue by FC (MeOH/CH₂Cl₂ 1:49 → 3:97 → 1:24 → 1:19 → 1:9 → 1:4) yielded the title compound (4.80 g, 55%). R_f = 0.50 (MeOH/CH₂Cl₂ 25 1:9). LC-MS: R_t = 4.99; ES-: 603.20.

Compounds of type W

30 (*rac.*)-(1*R*^{*}, 5*S*^{*})-7-(4-Hydroxyphenyl)-6-(methylphenethylcarbamoyl)-3,9-diazabicyclo[3.3.1]non-6-ene-3,9-dicarboxylic acid 3-*tert*-butyl ester 9-(2,2,2-trichloro-1,1-dimethylethyl) ester (**W1**)

To a susp. of bicyclononene **V1** (1.50 g, 2.66 mmol) in CHCl₃ (30 mL) was added methylphenethylamine (0.774 mL, 5.32 mmol), DMAP (32.5 mg, 0.266 mmol), HOBr (36 mg, 0.266 mmol), and EDC·HCl (765 mg, 3.99 mmol) were added successively. After 3 days at rt the mixture was diluted in CH₂Cl₂ and washed
5 with aq. 1M HCl (1x) and aq. sat. NaHCO₃ (1x). The org. extracts were dried over MgSO₄ and filtered. The solvents were removed under reduced pressure. Purification of the residue by FC (EtOAc/heptane 1:9 → 1:4 → 3:7 → 2:3 → 1:1 → 3:2 → 7:3) yielded the title compound as a colorless solid (0.93 g, 51%). R_f = 0.25 (EtOAc/heptane 1:1). LC-MS: R_t = 5.86; ES-: 678.14.

10

(*rac.*)-(1*R*^{*,} 5*S*^{*)}-7-[4-(2-Hydroxyethyl)phenyl]-6-(methylphenethylcarbamoyl)-3,9-diazabicyclo[3.3.1]non-6-ene-3,9-dicarboxylic acid 3-*tert*-butyl ester 9-(2,2,2-trichloro-1,1-dimethylethyl) ester (**W2**)

15 To a sol. of bicyclononene **V2** (7.09 g, 11.97 mmol) in CHCl₃ (140 mL) were added *N*-methylphenethylamine (3.48 mL, 24.0 mmol), DMAP (137 mg, 1.12 mmol), HOBr (151 mg, 1.12 mmol) and EDC·HCl (3.44 g, 18.0 mmol). The mixture was stirred at rt for 3 days, before it was diluted with CH₂Cl₂ and washed with aq. 1M HCl (1x) and aq. sat. NaHCO₃ (1x). The org. extracts were dried over MgSO₄ and filtered. The solvents were removed under reduced pressure. Purification of the residue purified by FC (MeOH/CH₂Cl₂ 1:49 → 3:122 → 4:121 → 1:24 → 1:9 → 1:4) yielded the title compound (3.46 g, 41%). R_f = 0.26 (MeOH/CH₂Cl₂ 1:19). LC-MS: R_t = 5.87; ES+: 708.40.

20 25 (*rac.*)-(1*R*^{*,} 5*S*^{*)}-7-[4-(3-Hydroxypropyl)phenyl]-6-(methylphenethylcarbamoyl)-3,9-diazabicyclo[3.3.1]non-6-ene-3,9-dicarboxylic acid 3-*tert*-butyl ester 9-(2,2,2-trichloro-1,1-dimethylethyl) ester (**W3**)

30 To a sol. of bicyclononene **V3** (7.59 g, 12.5 mmol) in CHCl₃ (150 mL) were added methylphenethylamine (3.63 mL, 25.0 mmol), DMAP (153 mg, 1.25 mmol), HOBr (169 mg, 1.25 mmol) and EDC·HCl (3.80 g, 19.2 mmol). The mixture was stirred at rt for 3 days before it was diluted in CH₂Cl₂ and washed

with aq. 1M HCl (1x) and aq. sat. NaHCO₃ (1x). The org. extracts were dried over MgSO₄ and filtered. The solvents were removed under reduced pressure. Purification of the residue by FC (MeOH/CH₂Cl₂ 1:49 → 3:97 → 1:24 → 1:9 → 1:4) yielded the title compound (3.18 g, 35%). R_f = 0.42 (MeOH/CH₂Cl₂ 1:19).

5 LC-MS: R_t = 5.99; ES+: 744.50.

(rac.)-(IR, SS*)-3-Acetyl-6-hydroxymethyl-7-{4-[3-(2-methoxybenzyloxy)propoxy]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-9-carboxylic acid 2,2,2-trichloro-1,1-dimethylethyl ester (X)*

10

To a sol. of bicyclononene S1 (2.29 g, 3.21 mmol) in CH₂Cl₂ (100 mL) was added BF₃Et₂O (0.460 mL, 3.66 mmol) at -78 °C. The mixture was stirred at -78 °C for 30 min and DIBAL (1M in toluene, 6.42 mL, 6.42 mmol) was added. After 75 min, ice was added and the mixture was allowed to warm up to rt. CH₂Cl₂ was 15 added and the mixture was washed with aq. 1M HCl (1x). The org. extracts were separated, dried over MgSO₄ and filtered. The solvents were removed under reduced pressure. Purification of the residue by FC (EtOAc/heptane 1:2 → 1:1 → EtOAc) yielded the title compound (1.01 g, 47%).

20 *(rac.)-(IR*, SS*)-3-Acetyl-6-formyl-7-{4-[3-(2-methoxybenzyloxy)propoxy]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-9-carboxylic acid 2,2,2-trichloro-1,1-dimethylethyl ester (Y)*

25 To a sol. of bicyclononene X (258 mg, 0.385 mmol) in CH₂Cl₂ (5 mL) was added to 0 °C Dess-Martin periodane (170 mg, 0.401 mmol) at 0 °C. After 45 min. at 0 °C a second portion periodane was added. The sol. was stirred for 15 min before the solvents were removed under reduced pressure. Direct purification of the residue by FC (EtOAc/heptane 2:3 → 1:1 → 3:2 → 7:3) yielded the title compound (188 mg, 73%). R_f = 0.49 (EtOAc). LC-MS: R_t = 6.18; ES+: 667.21.

30

(*rac.*)-(1*R*^{*}, 5*S*^{*})-3-Acetyl-7-{4-[3-(2-methoxybenzyloxy)propoxy]phenyl}-6-methylaminomethyl-3,9-diazabicyclo[3.3.1]non-6-ene-9-carboxylic acid 2,2,2-trichloro-1,1-dimethylethyl ester (Z)

5 To a sol. of bicyclononene Y (334 mg, 0.50 mmol) in MeOH (10 mL) methylamine (40% in water, 0.215 mL, 2.5 mmol) was added. The mixture was stirred at rt for 1 h and then cooled to 0 °C. NaBH₄ (20 mg, 0.50 mmol) was added. The mixture was stirred at rt for 4 h before K₂CO₃ (263 mg) was added. After evaporation under reduced pressure the residue was distributed between
10 EtOAc and water. The org. extracts were separated, dried over MgSO₄, and filtered. The solvents were removed under reduced pressure. Purification of the residue by RP18-MPLC yielded the title compound (130 mg, 38%). LC-MS: R_t = 1.00; ES+: 682.14.

15 **(*rac.*)-(1*R*^{*}, 5*S*^{*})-7-{4-[3-(2-Chlorophenoxy)propyl]phenyl}-3,9-diazabicyclo-[3.3.1]non-6-ene-3,6,9-tricarboxylic acid 3-*tert*-butyl ester 9-(2,2,2-trichloro-1,1-dimethylethyl) ester (AA)**

Bicyclononene H6 (1.71 g, 2.3 mmol) was dissolved in EtOH (50 mL). Aq. 1M
20 NaOH (50 mL) was added and the mixture was heated to 80 °C. The sol. was stirred for 5 h at 80 °C, then allowed to cool down to rt. After acidification to pH = 1-2 with aq. 1M HCl the mixture was extracted with EtOAc (3x). The combined org. extracts were dried over MgSO₄, filtered and the solvents were removed under reduced pressure. Purification of the residue by FC
25 (EtOAc/heptane 2:3 → 1:2 → 1:1) yielded the title compound (504 mg, 31 %). R_f = 0.30 (EtOAc/heptane 1:1). LC-MS: R_t = 6.21; ES-: 712.34.

30 **(*rac.*)-(1*R*^{*}, 5*S*^{*})-7-{4-[3-(2-Chlorophenoxy)propyl]phenyl}-6-{[2-(2-chlorophenyl)ethyl]methylcarbamoyl}-3,9-diazabicyclo[3.3.1]non-6-ene-3,9-dicarboxylic acid 3-*tert*-butyl ester 9-(2,2,2-trichloro-1,1-dimethylethyl) ester (AB)**

Bicyclononene AA (504 mg, 0.703 mmol) was dissolved in CHCl₃ (25 mL). [2-(2-chlorophenyl)ethyl]methylamine (Jaques B.; Wallace R. G., *Tetrahedron*, 1977, 33, 581; 238 mg, 1.40 mmol), DIPEA (0.240 mL, 1.40 mmol), DMAP (17 mg, 0.14 mmol), HOBr (19 mg, 0.10 mmol) and EDC·HCl (135 mg, 1.40 mmol) were added. The sol. was stirred at rt overnight. The mixture was diluted with CH₂Cl₂ and washed with water (1x). The org. extracts were dried over MgSO₄ and the solvents were removed under reduced pressure. Purification of the residue by filtration through silica gel yielded the title compound as a yellowish foam (336 mg, 55%).

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(*rac.*)-(IR*, 5S*)-7-{4-[3-(2-Chlorophenoxy)propyl]phenyl}-6-{{[2-(2-chlorophenyl)ethyl]methylcarbamoyl}-3,9-diazabicyclo[3.3.1]non-6-ene-9-carboxylic acid 2,2,2-trichloro-1,1-dimethylethyl ester hydrochloride salt (AC)

15 Bicyclononene AB (336 mg, 0.378 mmol) was dissolved in CH₂Cl₂ (3 mL). 4M HCl/dioxane (13 mL) was added and the mixture was stirred at rt for 2 h. The solvents were removed under reduced pressure. Drying the residue at high vacuum yielded the title compound as a colorless foam that was used without further purification. LC-MS: R_t = 5.26; ES+: 765.85.

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9-Methyl-7-oxo-3,9-diazabicyclo[3.3.1]nonane-3-carboxylic acid *tert*-butyl ester (AE)

25 (4-Benzyl-6-ethoxycarbonylmethyl-1-methylpiperazin-2-yl)acetic acid ethyl ester (Patent WO 92/05174) (71.0 g, 0.196 mol) was dissolved in MeOH (400 mL). TFA (77.8 mL, 1.02 mol) was added and the flask was purged with nitrogen. Pd/C (10%, 50% moisture, 3.6 g) was added. The flask was closed and purged with hydrogen (3x). After 1 day, the mixture was filtered through celite and washed with MeOH. The solvent was removed under reduced pressure and the 30 foamy residue (92.7 g) was dried under high vacuum. A sol. of tBuOK (117.2 g, 1.04 mol) in toluene (3.07 L) was heated to reflux. A sol. of the crude piperazine formerly obtained, dissolved in THF (300 mL), was added dropwise over 50 min.

The black mixture was stirred for 10 further min. and allowed to cool to rt. The mixture was cooled to 0 °C and AcOH (36.6 mL, 0.635 mol) was added. The solvents were removed under reduced pressure and the residue purified by FC (MeOH/CH₂Cl₂ 1:9 → 1:4 → 1:3). The fractions with an R_f-value close to 0.10
5 (MeOH/CH₂Cl₂ 1:9) were collected and the solvent removed under reduced pressure. The residue was dissolved in aq. 5M HCl (2 L) and the reaction mixture heated to reflux overnight. The mixture was allowed to cool to rt, then cooled to 0 °C with an ice bath. The pH was brought to 12 by adding carefully solid KOH. This mixture was extracted with CH₂Cl₂ (3x). The combined org. extracts were
10 dried over MgSO₄ and the solvents were removed under reduced pressure. The residue was suspended in CH₂Cl₂ (400 mL) and cooled to 0 °C. DIPEA (19.1 mL, 112 mmol) was added. A sol. of Boc₂O (24.3 g, 113 mmol) in CH₂Cl₂ (200 mL) was added dropwise. The mixture was stirred for 1 h at 0 °C, then 1 h at rt. The mixture was washed with aq. 10% Na₂CO₃ (2x). The org. extracts were dried
15 over MgSO₄, filtered and the solvents were evaporated under reduced pressure. The residue was purified by FC (EtOAc/heptane 1:1 → EtOAc). The title compound was obtained as a solid that could be recrystallized from heptane (15.6 g, 30%). R_f = 0.45 (MeOH/CH₂Cl₂ 1:9). LC-MS: R_t = 1.55; ES+: 254.16.

20 **(1*R*, 5*S*)-9-Methyl-7-oxo-3,9-diazabicyclo[3.3.1]nonane-3,6-dicarboxylic acid
3-*tert*-butyl ester 6-methyl ester (AF)**

To a susp. of (-)-bis[(S)-1-phenylethyl]amine hydrochloride (226 mg, 0.864 mmol) in THF (3 mL) at 0 °C was added dropwise *n*-BuLi (1.6M in hexane, 1.136 mL, 1.808 mmol). The mixture was stirred for 1 h at 0 °C, then cooled to -78 °C. A sol. of bicyclononane AE (200 mg, 0.786 mmol) in THF (2 mL) was added dropwise over 3 min. The reaction mixture was stirred for 3 h at -78 °C, then methylcyanoformate (0.081 mL, 1.02 mmol) was added. The reaction mixture was stirred for 30 min. at -78 °C and a sol. of AgNO₃ (191 mg, 1.124 mmol) in
25 H₂O/THF (1:1, 2 mL) was added. After 10 min., H₂O (1.5 mL) and AcOH (1.5 mL) were added and the reaction mixture was allowed to warm to rt. Ammoniac (25% in water) was added until the Ag-salt had completely dissolved. The
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reaction mixture was extracted with EtOAc (1x) and CH₂Cl₂ (2x). The combined org. extracts were dried over MgSO₄ and the solvents were removed under reduced pressure. Purification of the residue by FC (MeOH/CH₂Cl₂ 1:14) yielded the title compound (167 mg, 68%). R_f = 0.37 (MeOH/CH₂Cl₂ 1:9). LC-MS: R_t = 5 0.76; ES+: 313.10. ee = 82%.

(1*R*, 5*S*)-9-Methyl-7-oxo-3,9-diazabicyclo[3.3.1]nonane-3,6-dicarboxylic acid 6-benzyl ester 3-*tert*-butyl ester (AG)

10 To a susp. of (-)-bis[(S)-1-phenylethyl]amine hydrochloride (226 mg, 0.864 mmol) in THF (3 mL) at 0 °C was added dropwise *n*-BuLi (1.6M in hexane, 1.136 mL, 1.808 mmol). The mixture was stirred for 1 h at 0 °C, then cooled to -78 °C. A sol. of bicyclononane AE (200 mg, 0.786 mmol) in THF (2 mL) was added dropwise over 3 min. The reaction mixture was stirred for 3 h at -78 °C, then 15 methylcyanoformate (0.081 mL, 1.02 mmol) was added. The reaction mixture was stirred for 30 min. at -78 °C and a sol. of AgNO₃ (191 mg, 1.124 mmol) in H₂O/THF (1:1, 2 mL) was added. After 10 min., H₂O (1.5 mL) and AcOH (1.5 mL) were added and the reaction mixture was allowed to warm to rt. Ammoniac (25% in water) was added until the Ag-salt had completely dissolved. The 20 reaction mixture was extracted with EtOAc (1x) and CH₂Cl₂ (2x). The combined org. extracts were dried over MgSO₄ and the solvents were removed under reduced pressure. Purification of the residue by FC (MeOH/CH₂Cl₂ 1:14) yielded the title compound (150 mg, 49%). R_f = 0.50 (MeOH/CH₂Cl₂ 1:9). LC-MS: R_t = 0.87; ES+: 389.09. ee = 84%.

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(rac.)-(1*R, 5*S**)-3-Acetyl-7-{4-[2-(2-bromo-5-fluorophenoxy)ethyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6,9-dicarboxylic acid 6-ethyl ester 9-(2,2,2-trichloro-1,1-dimethylethyl) ester (AH)**

30 A sol. of bicyclononene S3 (900 mg, 1.60 mmol) in toluene (15 mL) was purged with N₂ (4x). 2-Bromo-5-fluorophenol (0.267 mL, 2.4 mmol), TMAD (344 mg, 2.00 mmol) and tributylphosphine (1.18 mL, 4.80 mmol) were added and the

reaction mixture was heated to reflux for 1 h. The mixture was allowed to cool to rt, and the solvents were removed under reduced pressure. Purification of the residue by FC (EtOAc/heptane 1:24 → 1:9 → 2:3 → 7:3) yielded the title compound (1.06 g, 90%). R_f = 0.58 (EtOAc). LC-MS: R_t = 6.52; ES+: 733.00.

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Compounds of type AJ

(*rac.*)-(IR*, 5S*)-3-Acetyl-7-{4-[2-(2-bromo-5-fluorophenoxy)ethyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6,9-dicarboxylic acid 9-(2,2,2-trichloro-1,1-dimethylethyl) ester (AJ1)

A mixture of bicyclononene AH (1.06 g, 1.44 mmol) in EtOH (30 mL) and aq. 1M NaOH (30 mL) was stirred efficiently at 80 °C for 2.5 h. The mixture was allowed to cool to rt, acidified with aq. 1M HCl, and extracted with EtOAc (3x).
15 The combined org. phases were dried over MgSO₄, filtered, and the solvents were removed under reduced pressure. Purification of the residue by FC (EtOAc/heptane 1:1 → 3:1 → EtOAc → MeOH/EtOAc 1:19 → 1:9) yielded the title compound (845 mg, 83%). R_f = 0.10 (EtOAc). LC-MS: R_t = 5.78; ES-: 702.81.

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(*rac.*)-(IR*, 5S*)-3-Acetyl-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6,9-dicarboxylic acid 9-(2,2,2-trichloro-1,1-dimethylethyl) ester (AJ2)

25 A sol. of bicyclononene AR1 (296 mg, 0.39 mmol) in MeOH (10 mL) at 0 °C was purged with N₂ (3x). Pd/C (10%, cat. amount) was added and the mixture was purged with H₂ (4x). The mixture was stirred under H₂ for 2 h at 0 °C, and was filtered through *Celite*. The solvents were removed under reduced pressure and the residue was dried under high vacuum (200 mg, 80%). It was used without
30 further purification. LC-MS: R_t = 5.83.

(*rac.*)-(1*R*^{*, 5*S*})-3-Acetyl-7-{4-[3-(3-fluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6,9-dicarboxylic acid 9-(2,2,2-trichloro-1,1-dimethyl-ethyl) ester (AJ3)

5 As for compound AJ2, but from AR2 (205 mg, 0.25 mmol), Pd/C (cat. amount) and MeOH (10 mL). The crude material (100 mg, 56%) was used without further purification. LC-MS: R_t = 5.81.

(*rac.*)-(1*R*^{*, 5*S*})-3-Acetyl-7-{4-[3-(2-bromo-5-fluorophenoxy)propyl]phenyl}-10 3,9-diazabicyclo[3.3.1]non-6-ene-6,9-dicarboxylic acid 9-(2,2,2-trichloro-1,1-dimethylethyl) ester (AJ4)

As for compound AJ1, but from bicyclononene BL (1.55 g, 2.07 mmol), EtOH (55 mL) and aq. 1M NaOH (55 mL). Purification of the residue by FC (EtOAc/heptane 1:1 → EtOAc → MeOH/EtOAc 1:9) yielded the title compound (1.17 g, 78%). R_f = 0.20 (EtOAc). R_t = 6.02; ES+: 721.12.

Compounds of type AK

20 (*rac.*)-(1*R*^{*, 5*S*})-7-{4-[2-(2-Bromo-5-fluorophenoxy)ethyl]phenyl}-6-(methylphenethylcarbamoyl)-3,9-diazabicyclo[3.3.1]non-6-ene-3,9-dicarboxylic acid 3-*tert*-butyl ester 9-(2,2,2-trichloro-1,1-dimethylethyl) ester (AK1)

A sol. of bicyclononene W2 (1.68 g, 2.37 mmol) in toluene (50 mL) was purged with N₂ (4x). 2-Bromo-5-fluorophenol (0.403 mL, 3.56 mmol), azodicarboxylic dipiperidine (897 mg, 3.56 mmol) and tributyl phosphine (1.62 mL, 7.12 mmol) were added. The mixture was heated to reflux for 2 h. The mixture was then allowed to cool to rt and the solvents were removed under reduced pressure. Purification of the residue by FC (EtOAc/heptane 1:19 → 1:9 → 1:4 → 1:1 → 30 3:1) yielded the title compound (1.88 g, 90%). R_f = 0.80 (EtOAc). LC-MS: R_t = 7.30.

(*rac.*)-(1*R*^{*}, 5*S*^{*})-6-[(2-Chlorobenzyl)cyclopropylcarbamoyl]-7-{4-[3-(2,3,6-tri-fluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-3,9-dicarbo-xylic acid 3-*tert*-butyl ester 9-(2,2,2-trichloro-1,1-dimethylethyl) ester (AK2)

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A mixture of bicyclononene **AY1** (12.15 g, 16 mmol), (2-chlorobenzyl)-cyclopropylamine (9.08 g, 50 mmol), DIPEA (10.9 mL, 64 mmol), DMAP (488 mg, 4 mmol), HOBr (2.43 g, 18 mmol) and EDC-HCl (4.60 g, 24 mmol) in CH₂Cl₂ (250 mL) was stirred overnight. The mixture was diluted with CH₂Cl₂, 10 and washed with aq. 1M HCl (3x) and with aq. sat. NaHCO₃ (1x). The org. extracts were dried over MgSO₄, filtered, and the solvents were removed under reduced pressure. Purification of the residue by FC (EtOAc/heptane 1:5 → 1:3 → 1:2) yielded the title compound (9.10 g, 63%). LC-MS: R_t = 7.68.

15 **(*rac.*)-(1*R*^{*}, 5*S*^{*})-6-(Benzylcyclopropylcarbamoyl)-7-{4-[3-(2,3,6-trifluoro-phenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-3,9-dicarboxylic acid 3-*tert*-butyl ester 9-(2,2,2-trichloro-1,1-dimethylethyl) ester (AK3)**

20 As for bicyclononene **AK2**, but from bicyclononene **AY1** (552 mg, 0.75 mmol), benzylcyclopropylamine (Loeppky, R. N.; *et al.*, *J. Org. Chem.*, **2000**, *65*, 96; 221 mg, 1.50 mmol), DIPEA (0.515 mL, 3.00 mmol), DMAP (23 mg, 0.19 mmol), HOBr (101 mg, 0.75 mmol) and EDC-HCl (216 mg, 1.12 mmol) in CH₂Cl₂ (7 mL). Purification by FC yielded the title compound (570 mg, 88%). LC-MS: R_t = 1.28.

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(*rac.*)-(1*R*^{*}, 5*S*^{*})-6-[(2-chlorobenzyl)ethylcarbamoyl]-7-{4-[3-(2,3,6-trifluoro-phenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-3,9-dicarboxylic acid 3-*tert*-butyl ester 9-(2,2,2-trichloro-1,1-dimethylethyl) ester (AK4)

30 As for bicyclononene **AK2**, but from bicyclononene **AY1** (552 mg, 0.75 mmol), (2-chlorobenzyl)ethylamine (Ishihara, Y; *et al.*; *Chem. Pharm. Bull.*, **1991**, *39*, 3225; 255 mg, 1.50 mmol), DIPEA (0.515 mL, 3.00 mmol), DMAP (23 mg, 0.19

mmol), HOBr (101 mg, 0.75 mmol) and EDC·HCl (216 mg, 1.12 mmol) in CH₂Cl₂ (7 mL). Purification by FC yielded the title compound (543 mg, 82%). LC-MS: R_t = 1.29.

5 (*rac.*)-(1*R*^{*}, 5*S*^{*})-6-[Cyclopropyl-(2-fluorobenzyl)carbamoyl]-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-3,9-dicarboxylic acid 3-*tert*-butyl ester 9-(2,2,2-trichloro-1,1-dimethylethyl) ester (AK5)

10 As for bicyclononene AK2, but from bicyclononene AY1 (552 mg, 0.75 mmol), cyclopropyl-(2-fluorobenzyl)amine (248 mg, 1.50 mmol), DIPEA (0.515 mL, 3.00 mmol), DMAP (23 mg, 0.19 mmol), HOBr (101 mg, 0.75 mmol) and EDC·HCl (216 mg, 1.12 mmol) in CH₂Cl₂ (7 mL). Purification by FC yielded the title compound (526 mg, 79%). LC-MS: R_t = 1.28.

15 (*rac.*)-(1*R*^{*}, 5*S*^{*})-6-[Cyclopropyl-(3-trifluoromethylbenzyl)carbamoyl]-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-3,9-dicarboxylic acid 3-*tert*-butyl ester 9-(2,2,2-trichloro-1,1-dimethylethyl) ester (AK6)

20 As for bicyclononene AK2, but from bicyclononene AY1 (552 mg, 0.75 mmol), cyclopropyl-(3-trifluoromethylbenzyl)amine (Brabander, H. J.; *et al.*; *J. Org. Chem.*, 1967, 32, 4053; 323 mg, 1.50 mmol), DIPEA (0.515 mL, 3.00 mmol), DMAP (23 mg, 0.19 mmol), HOBr (101 mg, 0.75 mmol) and EDC·HCl (216 mg, 1.12 mmol) in CH₂Cl₂ (7 mL). Purification by FC yielded the title compound (551 mg, 79%). LC-MS: R_t = 1.25.

30 (*rac.*)-(1*R*^{*}, 5*S*^{*})-6-[Cyclopropyl-(2-methylbenzyl)carbamoyl]-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-3,9-dicarboxylic acid 3-*tert*-butyl ester 9-(2,2,2-trichloro-1,1-dimethylethyl) ester (AK7)

As for bicyclononene **AK2**, but from bicyclononene **AY1** (552 mg, 0.75 mmol), cyclopropyl-(2-methylbenzyl)-amine (242 mg, 1.50 mmol), DIPEA (0.515 mL, 3.00 mmol), DMAP (23 mg, 0.19 mmol), HOBr (101 mg, 0.75 mmol) and EDC·HCl (216 mg, 1.12 mmol) in CH₂Cl₂ (7 mL). Purification by FC yielded the
5 title compound (553 mg, 84%). LC-MS: R_t = 1.29.

(*rac.*)-(1*R*^{*,} 5*S*^{*)}-6-{Cyclopropyl-[2-(4-methoxyphenoxy)ethyl]carbamoyl}-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-3,9-dicarboxylic acid 3-*tert*-butyl ester 9-(2,2,2-trichloro-1,1-dimethyl-
10 ethyl) ester (**AK8**)

As for bicyclononene **AK2**, but from bicyclononene **AY1** (552 mg, 0.75 mmol), cyclopropyl-[2-(4-methoxy-phenoxy)ethyl]amine (311 mg, 1.50 mmol), DIPEA (0.515 mL, 3.00 mmol), DMAP (23 mg, 0.19 mmol), HOBr (101 mg, 0.75 mmol)
15 and EDC·HCl (216 mg, 1.12 mmol) in CH₂Cl₂ (7 mL). Purification by FC yielded the title compound (566 mg, 82%). LC-MS: R_t = 1.28.

(*rac.*)-(1*R*^{*,} 5*S*^{*)}-6-{Cyclopropyl-[2-(2-methoxyphenoxy)ethyl]carbamoyl}-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-3,9-dicarboxylic acid 3-*tert*-butyl ester 9-(2,2,2-trichloro-1,1-dimethyl-
20 ethyl) ester (**AK9**)

As for bicyclononene **AK2**, but from bicyclononene **AY1** (552 mg, 0.75 mmol), cyclopropyl-[2-(2-methoxy-phenoxy)ethyl]amine (311 mg, 1.50 mmol), DIPEA (0.515 mL, 3.00 mmol), DMAP (23 mg, 0.19 mmol), HOBr (101 mg, 0.75 mmol)
25 and EDC·HCl (216 mg, 1.12 mmol) in CH₂Cl₂ (7 mL). Purification by FC yielded the title compound (570 mg, 82%). LC-MS: R_t = 1.28.

(*rac.*)-(1*R*^{*,} 5*S*^{*)}-6-[Cyclopropyl-(2-*m*-tolyloxyethyl)carbamoyl]-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-3,9-dicarboxylic acid 3-*tert*-butyl ester 9-(2,2,2-trichloro-1,1-dimethylethyl) ester
30 (**AK10**)

As for bicyclononene **AK2**, but from bicyclononene **AY1** (552 mg, 0.75 mmol), cyclopropyl-(2-*m*-tolyloxy-ethyl)amine (287 mg, 1.50 mmol), DIPEA (0.515 mL, 3.00 mmol), DMAP (23 mg, 0.19 mmol), HOBr (101 mg, 0.75 mmol) and EDC·HCl (216 mg, 1.12 mmol) in CH₂Cl₂ (7 mL). Purification by FC yielded the title compound (506 mg, 74%). LC-MS: R_t = 1.30.

(*rac*.)-(*1R*^{*,} *5S*^{*)}-6-{Cyclopropyl-[2-(3,4-dimethylphenoxy)ethyl]carbamoyl}-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-3,9-dicarboxylic acid 3-*tert*-butyl ester 9-(2,2,2-trichloro-1,1-dimethyl-ethyl) ester (AK11)

As for bicyclononene AK2, but from bicyclononene AY1 (552 mg, 0.75 mmol), cyclopropyl-[2-(3,4-dimethyl-phenoxy)ethyl]amine (462 mg, 1.50 mmol), DIPEA (0.515 mL, 3.00 mmol), DMAP (23 mg, 0.19 mmol), HOBt (101 mg, 0.75 mmol) and EDC·HCl (216 mg, 1.12 mmol) in CH₂Cl₂ (7 mL). Purification by FC yielded the title compound (693 mg, 100%). LC-MS: R_t = 1.28.

(*rac.*)-(IR*, 5S*)-6-(Cyclopropylphenethylcarbamoyl)-7-{4-[3-(2,3,6-trifluoro-
phenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-3,9-dicarboxylic
acid 3-*tert*-butyl ester 9-(2,2,2-trichloro-1,1-dimethylethyl) ester (AK12)

As for bicyclononene **AK2**, but from bicyclononene **AY1** (552 mg, 0.75 mmol), cyclopropylphenethylamine (Smith, P. W.; *et al.*; *J. Med. Chem.*, **1998**, *41*, 787; 242 mg, 1.50 mmol), DIPEA (0.515 mL, 3.00 mmol), DMAP (23 mg, 0.19 mmol), HOBr (101 mg, 0.75 mmol) and EDC·HCl (216 mg, 1.12 mmol) in CH₂Cl₂ (7 mL). Purification by FC yielded the title compound (510 mg, 77%). LC-MS: R_t = 1.28.

30 (rac.)-(1*R*^{*,} 5*S*^{*})-6-{{[2-(2-Chlorophenyl)ethyl]cyclopropylearbamoyl}-7-{{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-3,9-

**dicarboxylic acid 3-*tert*-butyl ester 9-(2,2,2-trichloro-1,1-dimethylethyl) ester
(AK13)**

As for bicyclononene AK2, but from bicyclononene AY1 (552 mg, 0.75 mmol),
5 [2-(2-chlorophenyl)ethyl]-cyclopropylamine (294 mg, 1.50 mmol), DIPEA (0.515 mL, 3.00 mmol), DMAP (23 mg, 0.19 mmol), HOBt (101 mg, 0.75 mmol) and EDC·HCl (216 mg, 1.12 mmol) in CH₂Cl₂ (7 mL). Purification by FC yielded the title compound (540 mg, 79%). LC-MS: R_t = 1.28.

10 **(*rac.*)-(1*R*^{*}, 5*S*^{*})-6-{Cyclopropyl-[2-(2,3-difluorophenyl)ethyl]carbamoyl}-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-3,9-dicarboxylic acid 3-*tert*-butyl ester 9-(2,2,2-trichloro-1,1-dimethylethyl) ester (AK14)**

15 As for bicyclononene AK2, but from bicyclononene AY1 (552 mg, 0.75 mmol), cyclopropyl-[2-(2,3-difluoro-phenyl)ethyl]amine (296 mg, 1.50 mmol), DIPEA (0.515 mL, 3.00 mmol), DMAP (23 mg, 0.19 mmol), HOBt (101 mg, 0.75 mmol) and EDC·HCl (216 mg, 1.12 mmol) in CH₂Cl₂ (7 mL). Purification by FC yielded the title compound (572 mg, 83%). LC-MS: R_t = 1.28.

20 **(*rac.*)-(1*R*^{*}, 5*S*^{*})-6-{Cyclopropyl-[2-(4-fluorophenyl)ethyl]carbamoyl}-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-3,9-dicarboxylic acid 3-*tert*-butyl ester 9-(2,2,2-trichloro-1,1-dimethylethyl) ester (AK15)**

25 As for bicyclononene AK2, but from bicyclononene AY1 (552 mg, 0.75 mmol), cyclopropyl-[2-(4-fluoro-phenyl)ethyl]amine (269 mg, 1.50 mmol), DIPEA (0.515 mL, 3.00 mmol), DMAP (23 mg, 0.19 mmol), HOBt (101 mg, 0.75 mmol) and EDC·HCl (216 mg, 1.12 mmol) in CH₂Cl₂ (7 mL). Purification by FC yielded the title compound (533 mg, 79%). LC-MS: R_t = 1.28.

(*rac.*)-(1*R*^{*}, 5*S*^{*})-6-[Cyclopropyl-(2-*o*-tolylethyl)carbamoyl]-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-3,9-dicarboxylic acid 3-*tert*-butyl ester 9-(2,2,2-trichloro-1,1-dimethylethyl) ester (AK16)

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As for bicyclononene AK2, but from bicyclononene AY1 (552 mg, 0.75 mmol), cyclopropyl-(2-*o*-tolylethyl)-amine (263 mg, 1.50 mmol), DIPEA (0.515 mL, 3.00 mmol), DMAP (23 mg, 0.19 mmol), HOBr (101 mg, 0.75 mmol) and EDC·HCl (216 mg, 1.12 mmol) in CH₂Cl₂ (7 mL). Purification by FC yielded the title compound (562 mg, 84%). LC-MS: R_t = 1.29.

(*rac.*)-(1*R*^{*}, 5*S*^{*})-6-[Cyclopropyl-(3,5-dimethoxybenzyl)carbamoyl]-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-3,9-dicarboxylic acid 3-*tert*-butyl ester 9-(2,2,2-trichloro-1,1-dimethylethyl) ester (AK17)

As for bicyclononene AK2, but from bicyclononene AY1 (552 mg, 0.75 mmol), cyclopropyl-(3,5-dimethoxybenzyl)amine (311 mg, 1.50 mmol), DIPEA (0.515 mL, 3.00 mmol), DMAP (23 mg, 0.19 mmol), HOBr (101 mg, 0.75 mmol) and EDC·HCl (216 mg, 1.12 mmol) in CH₂Cl₂ (7 mL). Purification by FC yielded the title compound (530 mg, 76%). LC-MS: R_t = 1.28.

(*rac.*)-(1*R*^{*}, 5*S*^{*})-6-[Cyclopropyl-(2-*p*-tolylethyl)carbamoyl]-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-3,9-dicarboxylic acid 3-*tert*-butyl ester 9-(2,2,2-trichloro-1,1-dimethylethyl) ester (AK18)

As for bicyclononene AK2, but from bicyclononene AY1 (552 mg, 0.75 mmol), cyclopropyl-(2-*p*-tolylethyl)-amine (263 mg, 1.50 mmol), DIPEA (0.515 mL, 3.00 mmol), DMAP (23 mg, 0.19 mmol), HOBr (101 mg, 0.75 mmol) and EDC·HCl (216 mg, 1.12 mmol) in CH₂Cl₂ (7 mL). Purification by FC yielded the title compound (530 mg, 79%). LC-MS: R_t = 1.30.

(*rac.*)-(1*R*^{*, 5S^{*})-6-{Cyclopropyl-[2-(2-hydroxyethyl)benzyl]carbamoyl}-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-3,9-dicarboxylic acid 3-*tert*-butyl ester 9-(2,2,2-trichloro-1,1-dimethylethyl)ester (AK19)}

As for bicyclononene AK2, but from bicyclononene AY1 (1.30, 1.77 mmol), (2-allylbenzyl)cyclopropyl-amine (992 mg, 5.30 mmol), DIPEA (1.01 mL, 7.08 mmol), DMAP (54 mg, 0.44 mmol), HOBr (263 mg, 1.95 mmol) and EDC·HCl (509 mg, 2.66 mmol) in CH₂Cl₂ (25 mL). Purification by FC yielded the intermediate compound (1.49 g, 93%). LC-MS: R_t = 7.81.

Then as for compound AT, but from the former intermediate compound (1.49 g, 1.65 mmol), NMO·H₂O (245 mg, 1.82 mmol), and OsO₄ (2.5% in *tert*-BuOH, 0.207 mL, 0.017 mmol) in THF (4 mL), *tert*-BuOH (2 mL) and water (1 mL). Purification of the residue by FC (EtOAc/heptane 1:4 → 2:3 → 3:2 → 4:1 → EtOAc) yielded the 2nd intermediate compound (866 mg, 56%). R_f = 0.50 (EtOAc). LC-MS: R_t = 6.95.

Then as for compound AU, but from the 2nd intermediate compound (866 mg, 0.922 mmol) and NaIO₄ (217 mg, 1.01 mmol) in THF (8 mL) and water (2 mL). Purification of the residue by FC (EtOAc/heptane 1:4 → 2:3) yielded the 3rd intermediate compound (751 mg, 90%). R_f = 0.75 (EtOAc). Finally as for compound AV, but from the 3rd intermediate compound (751 mg, 0.828 mmol) and NaBH₄ (35 mg, 0.9 mmol) in MeOH (10 mL). Purification of the residue by FC (EtOAc/heptane 2:3) yielded the title compound (599 mg, 80%). LC-MS: R_t = 7.30.

(*rac.*)-(1*R*^{*, 5S^{*})-7-{4-[3-(2-Bromo-5-fluorophenoxy)propyl]phenyl}-6-[2-chlorobenzyl]cyclopropylcarbamoyl]-3,9-diazabicyclo[3.3.1]non-6-ene-3,9-dicarboxylic acid 3-*tert*-butyl ester 9-(2,2,2-trichloro-1,1-dimethylethyl) ester (AK20)}

As for bicyclononene **AK2**, but from bicyclononene **AY2** (7.68 g, 9.86 mmol), (2-chlorobenzyl)cyclopropylamine (5.37, 29.6 mmol), DIPEA (6.75 mL, 39.4 mmol), DMAP (301 mg, 2.47 mmol), HOBr (1.46 mg, 10.8 mmol) and EDC·HCl (2.84 g, 14.8 mmol) in CH₂Cl₂ (150 mL). Purification by FC (EtOAc/heptane 1:4 → 3:7
5 → 2:3 → 1:1) yielded the title compound (3.7 g, 40%). R_f = 0.55 (EtOAc/heptane 1:1). LC-MS: R_t = 7.97.

(*rac.*)-(IR*, 5S*)-6-(Benzylcyclopropylcarbamoyl)-7-{4-[3-(2-bromo-5-fluoro-phenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-3,9-dicarboxylic
10 acid 3-*tert*-butyl ester 9-(2,2,2-trichloro-1,1-dimethylethyl) ester (**AK21**)

As for bicyclononene **AK2**, but from bicyclononene **AY2** (779 mg, 1.00 mmol), benzylcyclopropylamine (Loeppky, R. N.; *et al.*, *J. Org. Chem.*, 2000, 65, 96; 1.27 g, 1.50 mmol), DIPEA (0.684 mL, 4.00 mmol), DMAP (30 mg, 0.25 mmol),
15 HOBr (135 mg, 1.00 mmol) and EDC·HCl (287 mg, 1.50 mmol) in CH₂Cl₂ (10 mL). Purification by FC yielded the title compound (520 mg, 57%). LC-MS: R_t = 7.79.

(*rac.*)-(IR*, 5S*)-7-{4-[3-(2-Bromo-5-fluorophenoxy)propyl]phenyl}-6-[(2-chlorobenzyl)ethylcarbamoyl]-3,9-diazabicyclo[3.3.1]non-6-ene-3,9-dicarboxylic acid 3-*tert*-butyl ester 9-(2,2,2-trichloro-1,1-dimethylethyl) ester (**AK22**)

As for bicyclononene **AK2**, but from bicyclononene **AY2** (779 mg, 1.00 mmol), (2-chlorobenzyl)ethylamine (Ishihara, Y; *et al.*; *Chem. Pharm. Bull.*, 1991, 39, 25 3225; 254 mg, 1.50 mmol), DIPEA (0.684 mL, 4.00 mmol), DMAP (30 mg, 0.25 mmol), HOBr (135 mg, 1.00 mmol) and EDC·HCl (287 mg, 1.50 mmol) in CH₂Cl₂ (10 mL). Purification by FC yielded the title compound (475 mg, 51%). LC-MS: R_t = 7.82.

30 (*rac.*)-(IR*, 5S*)-7-{4-[3-(2-Bromo-5-fluorophenoxy)propyl]phenyl}-6-[cyclopropyl-(2-fluorobenzyl)carbamoyl]-3,9-diazabicyclo[3.3.1]non-6-ene-

3,9-dicarboxylic acid 3-*tert*-butyl ester 9-(2,2,2-trichloro-1,1-dimethylethyl)ester (AK23)

As for bicyclononene AK2, but from bicyclononene AY2 (779 mg, 1.00 mmol),
5 cyclopropyl-(2-fluorobenzyl)amine (247 mg, 1.50 mmol), DIPEA (0.684 mL, 4.00
mmol), DMAP (30 mg, 0.25 mmol), HOBr (135 mg, 1.00 mmol) and EDC·HCl
(287 mg, 1.50 mmol) in CH₂Cl₂ (10 mL). Purification by FC yielded the title
compound (465 mg, 50%). LC-MS: R_t = 7.69.

10 **(*rac.*)-(IR*, 5S*)-7-{4-[3-(2-Bromo-5-fluorophenoxy)propyl]phenyl}-6-[cyclo-
propyl-(3-trifluoromethylbenzyl)carbamoyl]-3,9-diazabicyclo[3.3.1]non-6-
ene-3,9-dicarboxylic acid 3-*tert*-butyl ester 9-(2,2,2-trichloro-1,1-dimethylethyl)
ester (AK24)**

15 As for bicyclononene AK2, but from bicyclononene AY2 (779 mg, 1.00 mmol),
cyclopropyl-(3-trifluoromethylbenzyl)amine (Brabander, H. J.; *et al.*; *J. Org.
Chem.*, 1967, 32, 4053; 323 mg, 1.50 mmol), DIPEA (0.684 mL, 4.00 mmol),
DMAP (30 mg, 0.25 mmol), HOBr (135 mg, 1.00 mmol) and EDC·HCl (287 mg,
1.50 mmol) in CH₂Cl₂ (10 mL). Purification by FC yielded the title compound
20 (345 mg, 35%). LC-MS: R_t = 7.76.

25 **(*rac.*)-(IR*, 5S*)-7-{4-[3-(2-Bromo-5-fluorophenoxy)propyl]phenyl}-6-[cyclopropyl-(2-methylbenzyl)carbamoyl]-3,9-diazabicyclo[3.3.1]non-6-ene-
3,9-dicarboxylic acid 3-*tert*-butyl ester 9-(2,2,2-trichloro-1,1-dimethylethyl)
ester (AK25)**

As for bicyclononene AK2, but from bicyclononene AY2 (779 mg, 1.00 mmol),
cyclopropyl-(2-methylbenzyl)amine (242 mg, 1.50 mmol), DIPEA (0.684 mL,
4.00 mmol), DMAP (30 mg, 0.25 mmol), HOBr (135 mg, 1.00 mmol) and
30 EDC·HCl (287 mg, 1.50 mmol) in CH₂Cl₂ (10 mL). Purification by FC yielded
the title compound (722 mg, 78%). LC-MS: R_t = 7.77.

(rac.)-(IR, 5S*)-7-{4-[3-(2-Bromo-5-fluorophenoxy)propyl]phenyl}-6-[cyclopropyl-[2-(4-methoxyphenoxy)ethyl]carbamoyl]-3,9-diazabicyclo[3.3.1]non-6-ene-3,9-dicarboxylic acid 3-*tert*-butyl ester 9-(2,2,2-trichloro-1,1-dimethyl-ethyl) ester (AK26)*

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As for bicyclononene AK2, but from bicyclononene AY2 (779 mg, 1.00 mmol), cyclopropyl-(4-methoxyphenoxy)methylamine (311 mg, 1.50 mmol), DIPEA (0.684 mL, 4.00 mmol), DMAP (30 mg, 0.25 mmol), HOBr (135 mg, 1.00 mmol) and EDC·HCl (287 mg, 1.50 mmol) in CH₂Cl₂ (10 mL). Purification by FC yielded the title compound (579 mg, 60%). LC-MS: R_t = 7.64.

(rac.)-(IR, 5S*)-7-{4-[3-(2-Bromo-5-fluorophenoxy)propyl]phenyl}-6-[cyclopropyl-(2-*m*-tolyloxyethyl)carbamoyl]-3,9-diazabicyclo[3.3.1]non-6-ene-3,9-dicarboxylic acid 3-*tert*-butyl ester 9-(2,2,2-trichloro-1,1-dimethylethyl) ester (AK27)*

As for bicyclononene AK2, but from bicyclononene AY2 (779 mg, 1.00 mmol), cyclopropyl-*m*-tolyloxymethylamine (311 mg, 1.50 mmol), DIPEA (0.684 mL, 4.00 mmol), DMAP (30 mg, 0.25 mmol), HOBr (135 mg, 1.00 mmol) and EDC·HCl (287 mg, 1.50 mmol) in CH₂Cl₂ (10 mL). Purification by FC yielded the title compound (340 mg, 36%). LC-MS: R_t = 7.83.

(rac.)-(IR, 5S*)-7-{4-[3-(2-Bromo-5-fluorophenoxy)propyl]phenyl}-6-[cyclopropyl-[2-(3,4-dimethylphenoxy)ethyl]carbamoyl]-3,9-diazabicyclo[3.3.1]-non-6-ene-3,9-dicarboxylic acid 3-*tert*-butyl ester 9-(2,2,2-trichloro-1,1-dimethylethyl) ester (AK28)*

As for bicyclononene AK2, but from bicyclononene AY2 (779 mg, 1.00 mmol), cyclopropyl-(3,4-dimethylphenoxy)methylamine (308 mg, 1.50 mmol), DIPEA (0.684 mL, 4.00 mmol), DMAP (30 mg, 0.25 mmol), HOBr (135 mg, 1.00 mmol) and EDC·HCl (287 mg, 1.50 mmol) in CH₂Cl₂ (10 mL). Purification by FC yielded the title compound (470 mg, 49%). LC-MS: R_t = 7.93.

(rac.)-(1R, 5S*)-7-{4-[3-(2-Bromo-5-fluorophenoxy)propyl]phenyl}-6-(cyclopropylphenethylcarbamoyl)-3,9-diazabicyclo[3.3.1]non-6-ene-3,9-dicarboxylic acid 3-*tert*-butyl ester 9-(2,2,2-trichloro-1,1-dimethylethyl) ester (AK29)*

5

As for bicyclononene AK2, but from bicyclononene AY2 (779 mg, 1.00 mmol), cyclopropyl-phenethylamine (Smith, P. W.; *et al.*; *J. Med. Chem.*, **1998**, *41*, 787; 242 mg, 1.50 mmol), DIPEA (0.684 mL, 4.00 mmol), DMAP (30 mg, 0.25 mmol), HOBr (135 mg, 1.00 mmol) and EDC·HCl (287 mg, 1.50 mmol) in CH₂Cl₂ (10 mL). Purification by FC yielded the title compound (449 mg, 49%). LC-MS: R_t = 7.72.

(rac.)-(1R, 5S*)-7-{4-[3-(2-Bromo-5-fluorophenoxy)propyl]phenyl}-6-{[2-(2-chlorophenyl)ethyl]cyclopropylcarbamoyl}-3,9-diazabicyclo[3.3.1]non-6-ene-3,9-dicarboxylic acid 3-*tert*-butyl ester 9-(2,2,2-trichloro-1,1-dimethylethyl) ester (AK30)*

As for bicyclononene AK2, but from bicyclononene AY2 (779 mg, 1.00 mmol), [2-(2-chloro-phenyl)ethyl]cyclopropylamine (294 mg, 1.50 mmol), DIPEA (0.684 mL, 4.00 mmol), DMAP (30 mg, 0.25 mmol), HOBr (135 mg, 1.00 mmol) and EDC·HCl (287 mg, 1.50 mmol) in CH₂Cl₂ (10 mL). Purification by FC yielded the title compound (605 mg, 63%). LC-MS: R_t = 7.89.

(rac.)-(1R, 5S*)-7-{4-[3-(2-Bromo-5-fluorophenoxy)propyl]phenyl}-6-{cyclopropyl-[2-(2,3-difluorophenyl)ethyl]carbamoyl}-3,9-diazabicyclo[3.3.1]non-6-ene-3,9-dicarboxylic acid 3-*tert*-butyl ester 9-(2,2,2-trichloro-1,1-dimethyl-ethyl) ester (AK31)*

As for bicyclononene AK2, but from bicyclononene AY2 (779 mg, 1.00 mmol), cyclopropyl-[2-(2,3-difluorophenyl)ethyl]amine (296 mg, 1.50 mmol), DIPEA (0.684 mL, 4.00 mmol), DMAP (30 mg, 0.25 mmol), HOBr (135 mg, 1.00 mmol)

and EDC·HCl (287 mg, 1.50 mmol) in CH₂Cl₂ (10 mL). Purification by FC yielded the title compound (670 mg, 70%). LC-MS: R_t = 7.70.

5 *(rac.)-(1R*, 5S*)-7-{4-[3-(2-Bromo-5-fluorophenoxy)propyl]phenyl}-6-[cyclo-*
propyl-[2-(4-fluorophenyl)ethyl]carbamoyl]-3,9-diazabicyclo[3.3.1]non-6-
ene-3,9-dicarboxylic acid 3-tert-butyl ester 9-(2,2,2-trichloro-1,1-dimethyl-
ethyl) ester (AK32)

10 As for bicyclononene AK2, but from bicyclononene AY2 (779 mg, 1.00 mmol),
cyclopropyl-[2-(4-fluorophenyl)ethyl]amine (269 mg, 1.50 mmol), DIPEA (0.684
mL, 4.00 mmol), DMAP (30 mg, 0.25 mmol), HOBr (135 mg, 1.00 mmol) and
EDC·HCl (287 mg, 1.50 mmol) in CH₂Cl₂ (10 mL). Purification by FC yielded
the title compound (638 mg, 68%). LC-MS: R_t = 7.70.

15 *15 (rac.)-(1R*, 5S*)-7-{4-[3-(2-Bromo-5-fluorophenoxy)propyl]phenyl}-6-[cyclo-*
propyl-(2-o-tolylethyl)carbamoyl]-3,9-diazabicyclo[3.3.1]non-6-ene-3,9-
dicarboxylic acid 3-tert-butyl ester 9-(2,2,2-trichloro-1,1-dimethylethyl) ester
(AK33)

20 As for bicyclononene AK2, but from bicyclononene AY2 (779 mg, 1.00 mmol),
cyclopropyl-(2-o-tolylethyl)amine (263 mg, 1.50 mmol), DIPEA (0.684 mL, 4.00
mmol), DMAP (30 mg, 0.25 mmol), HOBr (135 mg, 1.00 mmol) and EDC·HCl
(287 mg, 1.50 mmol) in CH₂Cl₂ (10 mL). Purification by FC yielded the title
compound (659 mg, 70%). LC-MS: R_t = 7.58.

25

1:1 Mixture of *(rac.)-(1R*, 5S*)-7-{4-[3-(2-bromo-5-fluorophenoxy)propyl]-*
phenyl}-6-[(2R)-2-hydroxy-2-phenylethyl)methylcarbamoyl]-3,9-diaza-*
bicyclo[3.3.1]non-6-ene-3,9-dicarboxylic acid 3-tert-butyl ester 9-(2,2,2-
trichloro-1,1-dimethylethyl) ester and *(rac.)-(1R*, 5S*)-7-{4-[3-(2-bromo-5-*
30 *fluorophenoxy)propyl]phenyl}-6-[(2S*)-2-hydroxy-2-phenylethyl)methyl-*
carbamoyl]-3,9-diazabicyclo[3.3.1]non-6-ene-3,9-dicarboxylic acid 3-tert-
butyl ester 9-(2,2,2-trichloro-1,1-dimethylethyl) ester (AK34)

As for bicyclononene **AK2**, but from bicyclononene **AY2** (779 mg, 1.00 mmol), (*rac.*)-2-methylamino-1-phenylethanol (310 mg, 1.50 mmol), DIPEA (0.684 mL, 4.00 mmol), DMAP (30 mg, 0.25 mmol), HOBr (135 mg, 1.00 mmol) and 5 EDC·HCl (287 mg, 1.50 mmol) in CH₂Cl₂ (10 mL). Purification by FC yielded the title compounds (456 mg, 50%). LC-MS: R_t = 7.42.

(*rac.*)-(1*R*^{*}, 5*S*^{*})-7-{4-[3-(2-Bromo-5-fluorophenoxy)propyl]phenyl}-6-[cyclopropyl-(3,5-dimethoxybenzyl)carbamoyl]-3,9-diazabicyclo[3.3.1]non-6-10 ene-3,9-dicarboxylic acid 3-*tert*-butyl ester 9-(2,2,2-trichloro-1,1-dimethyl-ethoxy) ester (**AK35**)

As for bicyclononene **AK2**, but from bicyclononene **AY2** (779 mg, 1.00 mmol), cyclopropyl-(3,5-dimethoxybenzyl)amine (311 mg, 1.50 mmol), DIPEA (0.684 15 mL, 4.00 mmol), DMAP (30 mg, 0.25 mmol), HOBr (135 mg, 1.00 mmol) and EDC·HCl (287 mg, 1.50 mmol) in CH₂Cl₂ (10 mL). Purification by FC yielded the title compound (736 mg, 76%). LC-MS: R_t = 7.73.

(*rac.*)-(1*R*^{*}, 5*S*^{*})-7-{4-[3-(2-Bromo-5-fluorophenoxy)propyl]phenyl}-6-[cyclo-20 propyl-(2-*p*-tolylethyl)carbamoyl]-3,9-diazabicyclo[3.3.1]non-6-ene-3,9-dicarboxylic acid 3-*tert*-butyl ester 9-(2,2,2-trichloro-1,1-dimethylethyl) ester (**AK36**)

As for bicyclononene **AK2**, but from bicyclononene **AY2** (779 mg, 1.00 mmol), 25 cyclopropyl-(2-*p*-tolylethyl)amine (263 mg, 1.50 mmol), DIPEA (0.684 mL, 4.00 mmol), DMAP (30 mg, 0.25 mmol), HOBr (135 mg, 1.00 mmol) and EDC·HCl (287 mg, 1.50 mmol) in CH₂Cl₂ (10 mL). Purification by FC yielded the title compound (718 mg, 77%). LC-MS: R_t = 7.73.

30 (*rac.*)-(1*R*^{*}, 5*S*^{*})-6-[(2-Allylbenzyl)cyclopropylcarbamoyl]-7-{4-[3-(2-bromo-5-fluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-3,9-

**dicarbo-xylic acid 3-*tert*-butyl ester 9-(2,2,2-trichloro-1,1-dimethylethyl) ester
(AK37)**

As for bicyclononene AK2, but from bicyclononene AY2 (1.45, 2.00 mmol), (2-allylbenzyl)cyclopropylamine (1.12 g, 6.00 mmol), DIPEA (1.37 mL, 8.00 mmol), DMAP (62 mg, 0.50 mmol), HOBr (298 mg, 2.20 mmol) and EDC·HCl (576 mg, 3.00 mmol) in CH₂Cl₂ (20 mL). Purification by FC yielded the intermediate compound (1.77 g, 93%). LC-MS: R_t = 7.95.

Then as for compound AT, but from the former intermediate compound (1.77 g, 1.86 mmol), NMO·H₂O (516 mg, 3.82 mmol), and OsO₄ (2.5% in *tert*-BuOH, 0.276 mL, 0.023 mmol) in THF (40 mL), *tert*-BuOH (20 mL) and water (10 mL). Purification of the residue by FC (EtOAc/heptane 1:4 → 2:3 → 3:2 → 4:1 → EtOAc) yielded the 2nd intermediate compound (548 mg, 27%). R_f = 0.60 (EtOAc). LC-MS: R_t = 7.43; ES+: 980.18.

15 Then as for compound AU, but from the 2nd intermediate compound (928 mg, 0.945 mmol) and NaIO₄ (222 mg, 1.04 mmol) in THF (8 mL) and water (2 mL). Purification of the residue by FC (EtOAc/heptane 1:4 → 2:3) yielded the 3rd intermediate compound (868 mg, 97%). R_f = 0.80 (EtOAc).

Finally as for compound AV, but from the 3rd intermediate compound (868 mg, 0.914 mmol) and NaBH₄ (38 mg, 1.0 mmol) in MeOH (10 mL). Purification of the residue by FC (EtOAc/heptane 2:3) yielded the title compound (603 mg, 69%). LC-MS: R_t = 7.44.

**(*rac.*)-(1*R*^{*}, 5*S*^{*})-6-[(2-Chlorobenzyl)cyclopropylcarbamoyl]-7-{4-[2-(2,3,5-trimethylphenoxy)ethyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-3,9-dicarboxylic acid 3-*tert*-butyl ester 9-(2,2,2-trichloro-1,1-dimethylethyl) ester
(AK38)**

As for bicyclononene AK2, but from bicyclononene AY3 (411 mg, 0.58 mmol), (2-chlorobenzyl)-cyclopropylamine (211 mg, 1.16 mmol), DIPEA (0.397 mL, 2.32 mmol), DMAP (18 mg, 0.15 mmol), HOBr (86 mg, 0.64 mmol) and

EDC·HCl (167 mg, 0.87 mmol) in CH₂Cl₂ (8 mL). Purification by FC yielded the title compound (312 mg, 62%). LC-MS: R_t = 7.66.

(*rac.*)-(1*R*^{*,}, 5*S*^{*})-6-(Benzylcyclopropylcarbamoyl)-7-{4-[2-(2,3,5-trimethyl-phenoxy)ethyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-3,9-dicarboxylic acid 3-*tert*-butyl ester 9-(2,2,2-trichloro-1,1-dimethylethyl) ester (AK39)

As for bicyclononene AK2, but from bicyclononene AY3 (411 mg, 0.58 mmol), benzyl-cyclopropylamine (Loeppky, R. N.; *et al.*, *J. Org. Chem.*, **2000**, *65*, 96; 10 171 mg, 1.16 mmol), DIPEA (0.397 mL, 2.32 mmol), DMAP (18 mg, 0.15 mmol), HOBr (86 mg, 0.64 mmol) and EDC·HCl (167 mg, 0.87 mmol) in CH₂Cl₂ (8 mL). Purification by FC yielded the title compound (340 mg, 70%). LC-MS: R_t = 8.12.

(15) (*rac.*)-(1*R*^{*,}, 5*S*^{*})-6-[(2-Chlorobenzyl)ethylcarbamoyl]-7-{4-[2-(2,3,5-trimethyl-phenoxy)ethyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-3,9-dicarboxylic acid 3-*tert*-butyl ester 9-(2,2,2-trichloro-1,1-dimethylethyl) ester (AK40)

20 As for bicyclononene AK2, but from bicyclononene AY3 (411 mg, 0.58 mmol), (2-chlorobenzyl)-ethylamine (Ishihara, Y; *et al.*; *Chem. Pharm. Bull.*, **1991**, *39*, 3225; 197 mg, 1.16 mmol), DIPEA (0.397 mL, 2.32 mmol), DMAP (18 mg, 0.15 mmol), HOBr (86 mg, 0.64 mmol) and EDC·HCl (167 mg, 0.87 mmol) in CH₂Cl₂ (8 mL). Purification by FC yielded the title compound (374 mg, 74%). LC-MS: 25 R_t = 8.30.

(*rac.*)-(1*R*^{*,}, 5*S*^{*})-6-[Cyclopropyl-(2-fluorobenzyl)carbamoyl]-7-{4-[2-(2,3,5-tri-methylphenoxy)ethyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-3,9dicarboxylic acid 3-*tert*-butyl ester 9-(2,2,2-trichloro-1,1-dimethylethyl) ester (AK41)

As for bicyclononene **AK2**, but from bicyclononene **AY3** (411 mg, 0.58 mmol), cyclopropyl-(2-fluorobenzyl)amine (192 mg, 1.16 mmol), DIPEA (0.397 mL, 2.32 mmol), DMAP (18 mg, 0.15 mmol), HOBr (86 mg, 0.64 mmol) and EDC·HCl (167 mg, 0.87 mmol) in CH₂Cl₂ (8 mL). Purification by FC yielded the title
5 compound (350 mg, 70%). LC-MS: R_t = 8.13.

(*rac.*)-(1*R*^{*,} 5*S*^{*)}-6-[Cyclopropyl-(3-trifluoromethylbenzyl)carbamoyl]-7-{4-[2-(2,3,5-trimethylphenoxy)ethyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-3,9-dicarboxylic acid 3-*tert*-butyl ester 9-(2,2,2-trichloro-1,1-dimethylethyl)
10 ester (**AK42**)

As for bicyclononene **AK2**, but from bicyclononene **AY3** (411 mg, 0.58 mmol), cyclopropyl-(3-trifluoromethylbenzyl)amine (Brabander, H. J.; *et al.*; *J. Org. Chem.*, 1967, 32, 4053; 250 mg, 1.16 mmol), DIPEA (0.397 mL, 2.32 mmol),
15 DMAP (18 mg, 0.15 mmol), HOBr (86 mg, 0.64 mmol) and EDC·HCl (167 mg, 0.87 mmol) in CH₂Cl₂ (8 mL). Purification by FC yielded the title compound (294 mg, 56%). LC-MS: R_t = 8.16.

(*rac.*)-(1*R*^{*,} 5*S*^{*)}-6-[Cyclopropyl-(2-methylbenzyl)carbamoyl]-7-{4-[2-(2,3,5-trimethylphenoxy)ethyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-3,9-dicarboxylic acid 3-*tert*-butyl ester 9-(2,2,2-trichloro-1,1-dimethylethyl) ester
20 (**AK43**)

As for bicyclononene **AK2**, but from bicyclononene **AY3** (411 mg, 0.58 mmol),
25 cyclopropyl-(2-methylbenzyl)amine (187 mg, 1.16 mmol), DIPEA (0.397 mL, 2.32 mmol), DMAP (18 mg, 0.15 mmol), HOBr (86 mg, 0.64 mmol) and EDC·HCl (167 mg, 0.87 mmol) in CH₂Cl₂ (8 mL). Purification by FC yielded the title compound (294 mg, 56%). LC-MS: R_t = 8.15.

30 (*rac.*)-(1*R*^{*,} 5*S*^{*)}-6-{Cyclopropyl-[2-(4-methoxyphenoxy)ethyl]carbamoyl}-7-{4-[2-(2,3,5-trimethylphenoxy)ethyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-

ene-3,9-dicarboxylic acid 3-*tert*-butyl ester 9-(2,2,2-trichloro-1,1-dimethyl-ethyl) ester (AK44)

As for bicyclononene AK2, but from bicyclononene AY3 (411 mg, 0.58 mmol),
5 cyclopropyl-[2-(4-methoxyphenoxy)ethyl]amine (187 mg, 1.16 mmol), DIPEA
(0.397 mL, 2.32 mmol), DMAP (18 mg, 0.15 mmol), HOBt (86 mg, 0.64 mmol)
and EDC·HCl (167 mg, 0.87 mmol) in CH₂Cl₂ (8 mL). Purification by FC yielded
the title compound (159 mg, 30%). LC-MS: R_t = 7.93.

10 **(*rac.*)-(1*R*^{*,} 5*S*^{*})-6-{Cyclopropyl-[2-(3-methoxyphenoxy)ethyl]carbamoyl}-7-{4-[2-(2,3,5-trimethylphenoxy)ethyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-3,9-dicarboxylic acid 3-*tert*-butyl ester 9-(2,2,2-trichloro-1,1-dimethyl-ethyl) ester (AK45)**

15 As for bicyclononene AK2, but from bicyclononene AY3 (411 mg, 0.58 mmol),
cyclopropyl-[2-(3-methoxyphenoxy)ethyl]amine (287 mg, 1.16 mmol), DIPEA
(0.397 mL, 2.32 mmol), DMAP (18 mg, 0.15 mmol), HOBt (86 mg, 0.64 mmol)
and EDC·HCl (167 mg, 0.87 mmol) in CH₂Cl₂ (8 mL). Purification by FC yielded
the title compound (237 mg, 45%). LC-MS: R_t = 7.69.

20 **(*rac.*)-(1*R*^{*,} 5*S*^{*})-6-[Cyclopropyl-(2-*m*-tolyloxyethyl)carbamoyl]-7-{4-[2-(2,3,5-trimethylphenoxy)ethyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-3,9-dicarboxylic acid 3-*tert*-butyl ester 9-(2,2,2-trichloro-1,1-dimethylethyl) ester (AK46)**

25 As for bicyclononene AK2, but from bicyclononene AY3 (411 mg, 0.58 mmol),
cyclopropyl-(2-*m*-tolyloxyethyl)amine (222 mg, 1.16 mmol), DIPEA (0.397 mL,
2.32 mmol), DMAP (18 mg, 0.15 mmol), HOBt (86 mg, 0.64 mmol) and
EDC·HCl (167 mg, 0.87 mmol) in CH₂Cl₂ (8 mL). Purification by FC yielded the
30 title compound (185 mg, 36%). LC-MS: R_t = 8.12.

(*rac.*)-(1*R*^{*,}, 5*S*^{*})-6-(Cyclopropylphenethylcarbamoyl)-7-{4-[2-(2,3,6-trimethyl-phenoxy)ethyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-3,9-dicarboxylic acid 3-*tert*-butyl ester 9-(2,2,2-trichloro-1,1-dimethylethyl) ester (AK47)

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As for bicyclononene AK2, but from bicyclononene AY3 (590 mg, 0.83 mmol), cyclopropyl-phenethylamine (Smith, P. W.; *et al.*; *J. Med. Chem.*, 1998, 41, 787; 402 mg, 2.49 mmol), DIPEA (0.568 mL, 3.32 mmol), DMAP (25 mg, 0.21 mmol), HOBr (169 mg, 1.25 mmol) and EDC·HCl (239 mg, 1.25 mmol) in CH₂Cl₂ (10 mL). Purification by FC yielded the title compound (309 mg, 44%). LC-MS: R_t = 8.01.

(*rac.*)-(1*R*^{*,}, 5*S*^{*})-6-{[2-(2-Chlorophenyl)ethyl]cyclopropylcarbamoyl}-7-{4-[2-(2,3,6-trimethylphenoxy)ethyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-3,9-dicarboxylic acid 3-*tert*-butyl ester 9-(2,2,2-trichloro-1,1-dimethylethyl) ester (AK48)

As for bicyclononene AK2, but from bicyclononene AY3 (590 mg, 0.83 mmol), [2-(2-chloro-phenyl)ethyl]cyclopropylamine (487 mg, 2.49 mmol), DIPEA (0.568 mL, 3.32 mmol), DMAP (25 mg, 0.21 mmol), HOBr (169 mg, 1.25 mmol) and EDC·HCl (239 mg, 1.25 mmol) in CH₂Cl₂ (10 mL). Purification by FC yielded the title compound (272 mg, 37%). LC-MS: R_t = 8.20.

(*rac.*)-(1*R*^{*,}, 5*S*^{*})-6-{Cyclopropyl-[2-(2,3-difluorophenyl)ethyl]carbamoyl}-7-{4-[2-(2,3,6-trimethylphenoxy)ethyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-3,9-dicarboxylic acid 3-*tert*-butyl ester 9-(2,2,2-trichloro-1,1-dimethylethyl) ester (AK49)

As for bicyclononene AK2, but from bicyclononene AY3 (590 mg, 0.83 mmol), cyclopropyl-[2-(2,3-difluorophenyl)ethyl]amine (491 mg, 2.49 mmol), DIPEA (0.568 mL, 3.32 mmol), DMAP (25 mg, 0.21 mmol), HOBr (169 mg, 1.25 mmol)

and EDC·HCl (239 mg, 1.25 mmol) in CH₂Cl₂ (10 mL). Purification by FC yielded the title compound (309 mg, 42%). LC-MS: R_t = 7.98.

(rac.)-(1R, 5S*)-6-{Cyclopropyl-[2-(4-fluorophenyl)ethyl]carbamoyl}-7-{4-*

5 *[2-(2,3,6-trimethylphenoxy)ethyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-*

3,9-dicarboxylic acid 3-tert-butyl ester 9-(2,2,2-trichloro-1,1-dimethylethyl)

ester (AK50)

As for bicyclononene AK2, but from bicyclononene AY3 (590 mg, 0.83 mmol),
10 cyclopropyl-[2-(4-fluorophenyl)ethyl]amine (491 mg, 2.49 mmol), DIPEA (0.568
mL, 3.32 mmol), DMAP (25 mg, 0.21 mmol), HOEt (169 mg, 1.25 mmol) and
EDC·HCl (239 mg, 1.25 mmol) in CH₂Cl₂ (10 mL). Purification by FC yielded
the title compound (294 mg, 41%). LC-MS: R_t = 7.93.

15 *(rac.)-(1R*, 5S*)-6-[Cyclopropyl-(2-o-tolylethyl)carbamoyl]-7-{4-[2-(2,3,6-*
trimethylphenoxy)ethyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-3,9-dicar-
boxylic acid 3-tert-butyl ester 9-(2,2,2-trichloro-1,1-dimethylethyl) ester
(AK51)

20 As for bicyclononene AK2, but from bicyclononene AY3 (590 mg, 0.83 mmol),
cyclopropyl-(2-o-tolylethyl)amine (491 mg, 2.49 mmol), DIPEA (0.568 mL, 3.32
mmol), DMAP (25 mg, 0.21 mmol), HOEt (169 mg, 1.25 mmol) and EDC·HCl
(239 mg, 1.25 mmol) in CH₂Cl₂ (10 mL). Purification by FC yielded the title
compound (258 mg, 36%). LC-MS: R_t = 8.16.

25

1:1-Mixture of (rac.)-1R*, 5S*)-6-[(2R*)-2-hydroxy-2-phenylethyl)methyl-

carbamoyl]-7-{4-[2-(2,3,6-trimethylphenoxy)ethyl]phenyl}-3,9-diazabicyclo-

[3.3.1]non-6-ene-3,9-dicarboxylic acid 3-tert-butyl ester 9-(2,2,2-trichloro-1,1-

dimethylethyl) ester and (rac.)-(1R*, 5S*)-6-[(2S*)-2-hydroxy-2-

30 **phenylethyl)methylcarbamoyl]-7-{4-[2-(2,3,6-trimethylphenoxy)ethyl]-**
phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-3,9-dicarboxylic acid 3-tert-butyl
ester 9-(2,2,2-trichloro-1,1-dimethylethyl) ester (AK52)

As for bicyclononene AK2, but from bicyclononene AY3 (590 mg, 0.83 mmol), (rac.)-2-methylamino-1-phenylethanol (377 mg, 2.49 mmol), DIPEA (0.568 mL, 3.32 mmol), DMAP (25 mg, 0.21 mmol), HOBr (169 mg, 1.25 mmol) and 5 EDC·HCl (239 mg, 1.25 mmol) in CH₂Cl₂ (10 mL). Purification by FC yielded the title compounds (117 mg, 17%). LC-MS: R_t = 7.50.

(rac.)-(1*R*^{*,} 5*S*^{*})-6-[Cyclopropyl-(3,5-dimethoxybenzyl)carbamoyl]-7-{4-[2-(2,3,6-trimethylphenoxy)ethyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-3,9-10 dicarboxylic acid 3-*tert*-butyl ester 9-(2,2,2-trichloro-1,1-dimethylethyl) ester (AK53)

As for bicyclononene AK2, but from bicyclononene AY3 (590 mg, 0.83 mmol), cyclopropyl-(3,5-dimethoxybenzyl)amine (516 mg, 2.49 mmol), DIPEA (0.568 15 mL, 3.32 mmol), DMAP (25 mg, 0.21 mmol), HOBr (169 mg; 1.25 mmol) and EDC·HCl (239 mg, 1.25 mmol) in CH₂Cl₂ (10 mL). Purification by FC yielded the title compound (258 mg, 35%). LC-MS: R_t = 7.80.

(rac.)-(1*R*^{*,} 5*S*^{*})-6-[Cyclopropyl-(2-*p*-tolylethyl)carbamoyl]-7-{4-[2-(2,3,6-20 trimethylphenoxy)ethyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-3,9- dicarboxylic acid 3-*tert*-butyl ester 9-(2,2,2-trichloro-1,1-dimethylethyl) ester (AK54)

As for bicyclononene AK2, but from bicyclononene AY3 (590 mg, 0.83 mmol), 25 cyclopropyl-(2-*p*-tolylethyl)amine (426 mg, 2.49 mmol), DIPEA (0.568 mL, 3.32 mmol), DMAP (25 mg, 0.21 mmol), HOBr (169 mg, 1.25 mmol) and EDC·HCl (239 mg, 1.25 mmol) in CH₂Cl₂ (10 mL). Purification by FC yielded the title compound (235 mg, 32%). LC-MS: R_t = 8.16.

30 (rac.)-(1*R*^{*,} 5*S*^{*})-6-[Cyclopropyl-[2-(2-hydroxyethyl)benzyl]carbamoyl]-7-{4-[2-(2,3,6-trimethylphenoxy)ethyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-

3,9-dicarboxylic acid 3-*tert*-butyl ester 9-(2,2,2-trichloro-1,1-dimethylethyl)ester (AK55)

As for bicyclononene AK2, but from bicyclononene AY3 (1.18 g, 1.66 mmol),
5 (2-allylbenzyl)cyclopropylamine (932 mg, 4.98 mmol), DIPEA (1.14 mL, 6.64
mmol), DMAP (51 mg, 0.42 mmol), HOBr (338 mg, 2.50 mmol) and EDC·HCl
(478 mg, 2.50 mmol) in CH₂Cl₂ (20 mL). Purification by FC yielded the
intermediate compound (613 mg, 42%). LC-MS: R_t = 8.16.

Then as for compound AT, but from the former intermediate compound (613 mg,
10 0.697 mmol), NMO·H₂O (141 mg, 1.05 mmol), and OsO₄ (2.5% in *tert*-BuOH,
0.175 mL, 0.014 mmol) in THF (8 mL), *tert*-BuOH (4 mL) and water (2 mL).
Purification of the residue by FC (EtOAc/heptane 1:1 → EtOAc) yielded the 2nd
intermediate compound (348 mg, 55%). R_f = 0.05 (EtOAc/heptane 1:1). LC-MS:
R_t = 7.31.

15 Then as for compound AU, but from the 2nd intermediate compound (348 mg,
0.381 mmol) and NaIO₄ (122 mg, 0.571 mmol) in THF (6 mL) and water (2 mL).
Drying the residue under high vacuum yielded the 3rd intermediate compound
(269 mg, 80%) that was used without further purification. LC-MS: R_t = 7.29.

Finally as for compound AV, but from the 3rd intermediate compound (269 mg,
20 0.305 mmol) and NaBH₄ (13 mg, 0.34 mmol) in MeOH (5 mL). Purification of
the residue by FC (EtOAc/heptane 1:4 → 2:3 → 3:2 → 4:1) yielded the title
compound (210 mg, 78%). LC-MS: R_t = 7.55.

**(*rac.*)-(1*R*^{*}, 5*S*^{*})-6-[(2-Chlorobenzyl)cyclopropylcarbamoyl]-7-{4-[2-(2-
25 chloro-4,5-dimethylphenoxy)ethoxy]phenyl}-3,9-diazabicyclo[3.3.1]non-6-
ene-3,9-dicarboxylic acid 3-*tert*-butyl ester 9-(2,2,2-trichloro-1,1-
dimethylethyl)ester (AK56)**

As for bicyclononene AK2, but from bicyclononene AY4 (596 mg, 0.80 mmol),
30 (2-chlorobenzyl)-cyclopropylamine (290 mg, 1.60 mmol), DIPEA (0.548 mL, 3.2
mmol), DMAP (25 mg, 0.21 mmol), HOBr (135 mg, 1.00 mmol) and EDC·HCl

(307 mg, 1.55 mmol) in CH₂Cl₂ (5 mL). Purification by FC yielded the title compound (451 mg, 74%). LC-MS: R_t = 1.30.

5 *(rac.)-(IR*, 5S*)-6-[(2-Chlorobenzyl)ethylcarbamoyl]-7-{4-[2-(2-chloro-4,5-dimethylphenoxy)ethoxy]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-3,9-dicarboxylic acid 3-tert-butyl ester 9-(2,2,2-trichloro-1,1-dimethylethyl) ester (AK57)*

10 As for bicyclononene AK2, but from bicyclononene AY4 (596 mg, 0.80 mmol), (2-chlorobenzyl)-ethylamine (Ishihara, Y; *et al.*; *Chem. Pharm. Bull.*, 1991, 39, 3225; 290 mg, 1.60 mmol), DIPEA (0.548 mL, 3.2 mmol), DMAP (25 mg, 0.21 mmol), HOBt (135 mg, 1.00 mmol) and EDC·HCl (307 mg, 1.55 mmol) in CH₂Cl₂ (5 mL). Purification by FC yielded the title compound (596 mg, 83%). LC-MS: R_t = 1.30.

15 *(rac.)-(IR*, 5S*)-7-{4-[2-(2-Chloro-4,5-dimethylphenoxy)ethoxy]phenyl}-6-[cyclopropyl-(2-fluorobenzyl)carbamoyl]-3,9-diazabicyclo[3.3.1]non-6-ene-3,9-dicarboxylic acid 3-tert-butyl ester 9-(2,2,2-trichloro-1,1-dimethylethyl) ester (AK58)*

20 As for bicyclononene AK2, but from bicyclononene AY4 (596 mg, 0.80 mmol), cyclopropyl-(2-fluorobenzyl)amine (264 mg, 1.60 mmol), DIPEA (0.548 mL, 3.2 mmol), DMAP (25 mg, 0.21 mmol), HOBt (135 mg, 1.00 mmol) and EDC·HCl (307 mg, 1.55 mmol) in CH₂Cl₂ (5 mL). Purification by FC yielded the title compound (519 mg, 73%). LC-MS: R_t = 1.29.

25 *(rac.)-(IR*, 5S*)-7-{4-[2-(2-Chloro-4,5-dimethylphenoxy)ethoxy]phenyl}-6-[cyclopropyl-(3-trifluoromethylbenzyl)carbamoyl]-3,9-diazabicyclo[3.3.1]-non-6-ene-3,9-dicarboxylic acid 3-tert-butyl ester 9-(2,2,2-trichloro-1,1-dimethylethyl) ester (AK59)*

As for bicyclononene AK2, but from bicyclononene AY4 (596 mg, 0.80 mmol), cyclopropyl-(3-trifluoromethylbenzyl)amine (Brabander, H. J.; *et al.*; *J. Org. Chem.*, 1967, 32, 4053; 344 mg, 1.60 mmol), DIPEA (0.548 mL, 3.2 mmol), DMAP (25 mg, 0.21 mmol), HOEt (135 mg, 1.00 mmol) and EDC·HCl (307 mg, 1.55 mmol) in CH₂Cl₂ (5 mL). Purification by FC yielded the title compound (584 mg, 78%). LC-MS: R_t = 1.30.

(*rac.*)-(1*R*^{*,} 5*S*^{*)}-7-{4-[2-(2-Chloro-4,5-dimethylphenoxy)ethoxy]phenyl}-6-[cyclopropyl-(2-methylbenzyl)carbamoyl]-3,9-diazabicyclo[3.3.1]non-6-ene-10 3,9-dicarboxylic acid 3-*tert*-butyl ester 9-(2,2,2-trichloro-1,1-dimethylethyl)ester (AK60)

As for bicyclononene AK2, but from bicyclononene AY4 (596 mg, 0.80 mmol), cyclopropyl-(2-methylbenzyl)amine (258 mg, 1.60 mmol), DIPEA (0.548 mL, 3.2 mmol), DMAP (25 mg, 0.21 mmol), HOEt (135 mg, 1.00 mmol) and EDC·HCl (307 mg, 1.55 mmol) in CH₂Cl₂ (5 mL). Purification by FC yielded the title compound (569 mg, 48%). LC-MS: R_t = 1.30.

(*rac.*)-(1*R*^{*,} 5*S*^{*)}-7-{4-[2-(2-Chloro-4,5-dimethylphenoxy)ethoxy]phenyl}-6-[cyclopropyl-[2-(4-methoxyphenoxy)ethyl]carbamoyl]-3,9-diazabicyclo[3.3.1]non-6-ene-3,9-dicarboxylic acid 3-*tert*-butyl ester 9-(2,2,2-trichloro-1,1-dimethylethyl)ester (AK61)

As for bicyclononene AK2, but from bicyclononene AY4 (596 mg, 0.80 mmol), cyclopropyl-[2-(4-methoxyphenoxy)ethyl]amine (332 mg, 1.60 mmol), DIPEA (0.548 mL, 3.2 mmol), DMAP (25 mg, 0.21 mmol), HOEt (135 mg, 1.00 mmol) and EDC·HCl (307 mg, 1.55 mmol) in CH₂Cl₂ (5 mL). Purification by FC yielded the title compound (591 mg, 79%). LC-MS: R_t = 1.28.

30 (*rac.*)-(1*R*^{*,} 5*S*^{*)}-7-{4-[2-(2-Chloro-4,5-dimethylphenoxy)ethoxy]phenyl}-6-[cyclopropyl-[2-(3-methoxyphenoxy)ethyl]carbamoyl]-3,9-diazabicyclo-

[3.3.1]non-6-ene-3,9-dicarboxylic acid 3-*tert*-butyl ester 9-(2,2,2-trichloro-1,1-dimethylethyl) ester (AK62)

As for bicyclononene AK2, but from bicyclononene AY4 (596 mg, 0.80 mmol),
5 cyclopropyl-[2-(3-methoxyphenoxy)ethyl]amine (332 mg, 1.60 mmol), DIPEA
(0.548 mL, 3.2 mmol), DMAP (25 mg, 0.21 mmol), HOBr (135 mg, 1.00 mmol)
and EDC·HCl (307 mg, 1.55 mmol) in CH₂Cl₂ (5 mL). Purification by FC yielded
the title compound (584 mg, 77%). LC-MS: R_t = 1.28.

10 **(*rac*.)-(1*R*^{*,} 5*S*^{*)}-7-{4-[2-(2-Chloro-4,5-dimethylphenoxy)ethoxy]phenyl}-6-[cyclopropyl-(2-*m*-tolyloxyethyl)carbamoyl]-3,9-diazabicyclo[3.3.1]non-6-ene-3,9-dicarboxylic acid 3-*tert*-butyl ester 9-(2,2,2-trichloro-1,1-dimethylethyl) ester (AK63)**

15 As for bicyclononene AK2, but from bicyclononene AY4 (596 mg, 0.80 mmol), cyclopropyl-(2-*p*-tolyloxyethyl)amine (306 mg, 1.60 mmol), DIPEA (0.548 mL, 3.2 mmol), DMAP (25 mg, 0.21 mmol), HOBr (135 mg, 1.00 mmol) and EDC·HCl (307 mg, 1.55 mmol) in CH₂Cl₂ (5 mL). Purification by FC yielded the title compound (525 mg, 71%). LC-MS: R_t = 1.30.

20 **(*rac*.)-(1*R*^{*,} 5*S*^{*)}-7-{4-[2-(2-Chloro-4,5-dimethylphenoxy)ethoxy]phenyl}-6-(cyclopropylphenethylcarbamoyl)-3,9-diazabicyclo[3.3.1]non-6-ene-3,9-dicarboxylic acid 3-*tert*-butyl ester 9-(2,2,2-trichloro-1,1-dimethylethyl) ester (AK64)**

25 As for bicyclononene AK2, but from bicyclononene AY4 (596 mg, 0.80 mmol), cyclopropyl-phenethylamine (Smith, P. W.; *et al.*; *J. Med. Chem.*, **1998**, *41*, 787; 258 mg, 1.60 mmol), DIPEA (0.548 mL, 3.2 mmol), DMAP (25 mg, 0.21 mmol), HOBr (135 mg, 1.00 mmol) and EDC·HCl (307 mg, 1.55 mmol) in CH₂Cl₂ (5 mL). Purification by FC yielded the title compound (360 mg, 50%). LC-MS: R_t = 1.30.

(*rac.*)-(1*R*^{*}, 5*S*^{*})-7-{4-[2-(2-Chloro-4,5-dimethylphenoxy)ethoxy]phenyl}-6-{{[2-(2-chlorophenyl)ethyl]cyclopropylcarbamoyl}-3,9-diazabicyclo[3.3.1]non-6-ene-3,9-dicarboxylic acid 3-*tert*-butyl ester 9-(2,2,2-trichloro-1,1-dimethyl-ethyl) ester (AK65)}

5

As for bicyclononene AK2, but from bicyclononene AY4 (596 mg, 0.80 mmol), [2-(2-chlorophenyl)ethyl]cyclopropylamine (313 mg, 1.60 mmol), DIPEA (0.548 mL, 3.2 mmol), DMAP (25 mg, 0.21 mmol), HOBr (135 mg, 1.00 mmol) and EDC·HCl (307 mg, 1.55 mmol) in CH₂Cl₂ (5 mL). Purification by FC yielded the 10 title compound (572 mg, 76%). LC-MS: R_t = 1.30.

(*rac.*)-(1*R*^{*}, 5*S*^{*})-7-{4-[2-(2-Chloro-4,5-dimethylphenoxy)ethoxy]phenyl}-6-{{cyclopropyl-[2-(2,3-difluorophenyl)ethyl]carbamoyl}-3,9-diazabicyclo[3.3.1]-non-6-ene-3,9-dicarboxylic acid 3-*tert*-butyl ester 9-(2,2,2-trichloro-1,1-dimethylethyl) ester (AK66)}

As for bicyclononene AK2, but from bicyclononene AY4 (596 mg, 0.80 mmol), cyclopropyl-[2-(2,3-difluorophenyl)ethyl]amine (316 mg, 1.60 mmol), DIPEA (0.548 mL, 3.2 mmol), DMAP (25 mg, 0.21 mmol), HOBr (135 mg, 1.00 mmol) 20 and EDC·HCl (307 mg, 1.55 mmol) in CH₂Cl₂ (5 mL). Purification by FC yielded the title compound (584 mg, 79%). LC-MS: R_t = 1.29.

(*rac.*)-(1*R*^{*}, 5*S*^{*})-7-{4-[2-(2-Chloro-4,5-dimethylphenoxy)ethoxy]phenyl}-6-{{cyclopropyl-[2-(4-fluorophenyl)ethyl]carbamoyl}-3,9-diazabicyclo[3.3.1]non-25 6-ene-3,9-dicarboxylic acid 3-*tert*-butyl ester 9-(2,2,2-trichloro-1,1-dimethyl-ethyl) ester (AK67)}

As for bicyclononene AK2, but from bicyclononene AY4 (596 mg, 0.80 mmol), cyclopropyl-[2-(4-fluorophenyl)ethyl]amine (287 mg, 1.60 mmol), DIPEA (0.548 mL, 3.2 mmol), DMAP (25 mg, 0.21 mmol), HOBr (135 mg, 1.00 mmol) 30 and EDC·HCl (307 mg, 1.55 mmol) in CH₂Cl₂ (5 mL). Purification by FC yielded the title compound (616 mg, 84%). LC-MS: R_t = 1.28.

(*rac.*)-(1*R*^{*}, 5*S*^{*})-7-{4-[2-(2-Chloro-4,5-dimethylphenoxy)ethoxy]phenyl}-6-[cyclopropyl-(2-*o*-tolylethyl)carbamoyl]-3,9-diazabicyclo[3.3.1]non-6-ene-3,9-dicarboxylic acid 3-*tert*-butyl ester 9-(2,2,2-trichloro-1,1-dimethylethyl) ester
5 (AK68)

As for bicyclononene AK2, but from bicyclononene AY4 (596 mg, 0.80 mmol), cyclopropyl-(2-*o*-tolylethyl)amine (280 mg, 1.60 mmol), DIPEA (0.548 mL, 3.2 mmol), DMAP (25 mg, 0.21 mmol), HOBr (135 mg, 1.00 mmol) and EDC·HCl 10 (307 mg, 1.55 mmol) in CH₂Cl₂ (5 mL). Purification by FC yielded the title compound (556 mg, 76%). LC-MS: R_t = 1.28.

1:1-Mixture of (*rac.*)-(1*R*^{*}, 5*S*^{*})-7-{4-[2-(2-chloro-4,5-dimethylphenoxy)ethoxy]phenyl}-6-[(2*R*^{*})-2-hydroxy-2-phenylethyl)methylcarbamoyl]-3,9-diazabicyclo[3.3.1]non-6-ene-3,9-dicarboxylic acid 3-*tert*-butyl ester 9-(2,2,2-trichloro-1,1-dimethylethyl) ester and (*rac.*)-(1*R*^{*}, 5*S*^{*})-7-{4-[2-(2-chloro-4,5-dimethylphenoxy)ethoxy]phenyl}-6-[(2*S*^{*})-2-hydroxy-2-phenylethyl)methylcarbamoyl]-3,9-diazabicyclo[3.3.1]non-6-ene-3,9-dicarboxylic acid 3-*tert*-butyl ester 9-(2,2,2-trichloro-1,1-dimethylethyl) ester (AK69)
15

20 As for bicyclononene AK2, but from bicyclononene AY4 (596 mg, 0.80 mmol), (*rac.*)-2-methylamino-1-phenylethanol (242 mg, 1.60 mmol), DIPEA (0.548 mL, 3.2 mmol), DMAP (25 mg, 0.21 mmol), HOBr (135 mg, 1.00 mmol) and EDC·HCl (307 mg; 1.55 mmol) in CH₂Cl₂ (5 mL). Purification by FC yielded the 25 title compounds (380 mg, 54%). LC-MS: R_t = 1.23.

(*rac.*)-(1*R*^{*}, 5*S*^{*})-7-{4-[2-(2-Chloro-4,5-dimethylphenoxy)ethoxy]phenyl}-6-[cyclopropyl-(3,5-dimethoxybenzyl)carbamoyl]-3,9-diazabicyclo[3.3.1]non-6-ene-3,9-dicarboxylic acid 3-*tert*-butyl ester 9-(2,2,2-trichloro-1,1-dimethyl-30 ethyl) ester (AK70)

As for bicyclononene AK2, but from bicyclononene AY4 (596 mg, 0.80 mmol), cyclopropyl-(3,5-dimethoxybenzyl)amine (332 mg, 1.60 mmol), DIPEA (0.548 mL, 3.2 mmol), DMAP (25 mg, 0.21 mmol), HOBr (135 mg, 1.00 mmol) and EDC·HCl (307 mg, 1.55 mmol) in CH₂Cl₂ (5 mL). Purification by FC yielded the title compound (619 mg, 83%). LC-MS: R_t = 1.28.

(*rac.*)-(1*R*^{*,} 5*S*^{*/})-7-{4-[2-(2-Chloro-4,5-dimethylphenoxy)ethoxy]phenyl}-6-{cyclopropyl-(2-*p*-tolylethyl)carbamoyl}-3,9-diazabicyclo[3.3.1]non-6-ene-3,9-dicarboxylic acid 3-*tert*-butyl ester 9-(2,2,2-trichloro-1,1-dimethylethyl) ester (AK71)

As for bicyclononene AK2, but from bicyclononene AY4 (596 mg, 0.80 mmol), cyclopropyl-(2-*p*-tolylethyl)amine (280 mg, 1.60 mmol), DIPEA (0.548 mL, 3.2 mmol), DMAP (25 mg, 0.21 mmol), HOBr (135 mg, 1.00 mmol) and EDC·HCl (307 mg, 1.55 mmol) in CH₂Cl₂ (5 mL). Purification by FC yielded the title compound (619 mg, 83%). LC-MS: R_t = 1.28.

(*rac.*)-(1*R*^{*,} 5*S*^{*/})-7-{4-[2-(2-Chloro-4,5-dimethylphenoxy)ethoxy]phenyl}-6-{cyclopropyl-[2-(2-hydroxyethyl)benzyl]carbamoyl}-3,9-diazabicyclo[3.3.1]-non-6-ene-3,9-dicarboxylic acid 3-*tert*-butyl ester 9-(2,2,2-trichloro-1,1-dimethyl-ethyl) ester (AK72)

As for bicyclononene AK2, but from bicyclononene AY3 (1.00 g, 1.34 mmol), (2-allylbenzyl)cyclopropylamine (752 mg, 4.02 mmol), DIPEA (0.918 mL, 5.26 mmol), DMAP (41 mg, 0.34 mmol), HOBr (199 mg, 1.47 mmol) and EDC·HCl (385 mg, 2.01 mmol) in CH₂Cl₂ (15 mL). Purification by FC yielded the intermediate compound (845 mg, 69%). R_f = 0.45 (EtOAc/heptane 1:1). LC-MS: R_t = 7.85.

Then as for compound AT, but from the former intermediate compound (845 mg, 0.926 mmol), NMO·H₂O (150 mg, 1.11 mmol), and OsO₄ (2.5% in *tert*-BuOH, 0.173 mL, 0.014 mmol) in THF (8 mL), *tert*-BuOH (4 mL) and water (2 mL). Purification of the residue by FC (EtOAc/heptane 1:1 → EtOAc → MeOH/EtOAc

1:9) yielded the 2nd intermediate compound (616 mg, 70%). R_f = 0.05 (EtOAc/heptane 1:1). LC-MS: R_t = 7.04.

Then as for compound AU, but from the 2nd intermediate compound (616 mg, 0.649 mmol) and NaIO₄ (208 mg, 0.973 mmol) in THF (6 mL) and water (2 mL).

5 Drying the residue under high vacuum yielded the 3rd intermediate compound (477 mg, 80%) that was used without further purification. LC-MS: R_t = 7.43.

Finally as for compound AV, but from the 3rd intermediate compound (477 mg, 0.520 mmol) and NaBH₄ (22 mg, 0.57 mmol) in MeOH (5 mL). Purification of the residue by FC (EtOAc/heptane 1:4 → 2:3 → 3:2 → 4:1) yielded the title 10 compound (210 mg, 78%). R_f = 0.10 (EtOAc/heptane 1:1). LC-MS: R_t = 7.26.

(rac.)-(1R, 5S*)-7-{4-[2-(4-Bromophenoxy)ethoxy]phenyl}-6-[(2-chlorobenzyl)cyclopropylcarbamoyl]-3,9-diazabicyclo[3.3.1]non-6-ene-3,9-dicarboxylic acid 3-tert-butyl ester 9-(2,2,2-trichloro-1,1-dimethylethyl) ester (AK73)*

15

As for bicyclononene AK2, but from bicyclononene AY5 (534 mg, 0.7 mmol), (2-chlorobenzyl)-cyclopropylamine (200 mg, 1.10 mmol), DIPEA (0.479 mL, 2.8 mmol), DMAP (21 mg, 0.18 mmol), HOBr (113 mg, 0.84 mmol) and EDC·HCl (211 mg, 1.1 mmol) in CH₂Cl₂ (7 mL). Purification by FC yielded the title 20 compound (263 mg, 39%). LC-MS: R_t = 1.28.

(rac.)-(1R, 5S*)-6-(Benzylcyclopropylcarbamoyl)-7-{4-[2-(4-bromophenoxy)ethoxy]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-3,9-dicarboxylic acid 3-tert-butyl ester 9-(2,2,2-trichloro-1,1-dimethylethyl) ester (AK74)*

25

As for bicyclononene AK2, but from bicyclononene AY5 (534 mg, 0.7 mmol), benzylcyclopropyl-amine (Loeppky, R. N.; *et al.*, *J. Org. Chem.*, 2000, 65, 96; 162 mg, 1.10 mmol), DIPEA (0.479 mL, 2.8 mmol), DMAP (21 mg, 0.18 mmol), HOBr (113 mg, 0.84 mmol) and EDC·HCl (211 mg, 1.1 mmol) in CH₂Cl₂ (7 mL). 30 Purification by FC yielded the title compound (263 mg, 39%). LC-MS: R_t = 1.26.

(*rac.*)-(1*R*^{*,}, 5*S*^{*})-7-{4-[2-(4-Bromophenoxy)ethoxy]phenyl}-6-[(2-chlorobenzyl)ethylcarbamoyl]-3,9-diazabicyclo[3.3.1]non-6-ene-3,9-dicarboxylic acid 3-*tert*-butyl ester 9-(2,2,2-trichloro-1,1-dimethylethyl) ester (AK75)

5 As for bicyclononene AK2, but from bicyclononene AY5 (534 mg, 0.7 mmol), (2-chlorobenzyl)-ethylamine (Ishihara, Y; *et al.*; *Chem. Pharm. Bull.*, 1991, 39, 3225; 187 mg, 1.10 mmol), DIPEA (0.479 mL, 2.8 mmol), DMAP (21 mg, 0.18 mmol), HOBr (113 mg, 0.84 mmol) and EDC-HCl (211 mg, 1.1 mmol) in CH₂Cl₂ (7 mL). Purification by FC yielded the title compound (204 mg, 32%). LC-MS: R_t = 1.28.

10 15 (*rac.*)-(1*R*^{*,} 5*S*^{*})-7-{4-[2-(4-Bromophenoxy)ethoxy]phenyl}-6-[cyclopropyl-(2-fluorobenzyl)carbamoyl]-3,9-diazabicyclo[3.3.1]non-6-ene-3,9-dicarboxylic acid 3-*tert*-butyl ester 9-(2,2,2-trichloro-1,1-dimethylethyl) ester (AK76)

As for bicyclononene AK2, but from bicyclononene AY5 (534 mg, 0.7 mmol), cyclopropyl-(2-fluorobenzyl)amine (182 mg, 1.10 mmol), DIPEA (0.479 mL, 2.8 mmol), DMAP (21 mg, 0.18 mmol), HOBr (113 mg, 0.84 mmol) and EDC-HCl (211 mg, 1.1 mmol) in CH₂Cl₂ (7 mL). Purification by FC yielded the title compound (233 mg, 37%). LC-MS: R_t = 1.27.

20 25 (*rac.*)-(1*R*^{*,} 5*S*^{*})-7-{4-[2-(4-Bromophenoxy)ethoxy]phenyl}-6-[cyclopropyl-(3-trifluoromethylbenzyl)carbamoyl]-3,9-diazabicyclo[3.3.1]non-6-ene-3,9-dicarboxylic acid 3-*tert*-butyl ester 9-(2,2,2-trichloro-1,1-dimethylethyl) ester (AK77)

As for bicyclononene AK2, but from bicyclononene AY5 (534 mg, 0.7 mmol), cyclopropyl-(3-trifluoromethylbenzyl)amine (Brabander, H. J.; *et al.*; *J. Org. Chem.*, 1967, 32, 4053; 237 mg, 1.10 mmol), DIPEA (0.479 mL, 2.8 mmol), DMAP (21 mg, 0.18 mmol), HOBr (113 mg, 0.84 mmol) and EDC-HCl (211 mg, 1.1 mmol) in CH₂Cl₂ (7 mL). Purification by FC yielded the title compound (276 mg, 41%). LC-MS: R_t = 1.27.

(rac.)-(1R, 5S*)-7-{4-[2-(4-Bromophenoxy)ethoxy]phenyl}-6-[cyclopropyl-(2-methylbenzyl)carbamoyl]-3,9-diazabicyclo[3.3.1]non-6-ene-3,9-dicarboxylic acid 3-*tert*-butyl ester 9-(2,2,2-trichloro-1,1-dimethylethyl) ester (AK78)*

5

As for bicyclononene AK2, but from bicyclononene AY5 (534 mg, 0.7 mmol), cyclopropyl-(2-methylbenzyl)amine (178 mg, 1.10 mmol), DIPEA (0.479 mL, 2.8 mmol), DMAP (21 mg, 0.18 mmol), HOBt (113 mg, 0.84 mmol) and EDC·HCl (211 mg, 1.1 mmol) in CH₂Cl₂ (7 mL). Purification by FC yielded the title compound (171 mg, 27%). LC-MS: R_t = 1.26.

(rac.)-(1R, 5S*)-7-{4-[2-(4-Bromophenoxy)ethoxy]phenyl}-6-[cyclopropyl-[2-(4-methoxyphenoxy)ethyl]carbamoyl]-3,9-diazabicyclo[3.3.1]non-6-ene-3,9-dicarboxylic acid 3-*tert*-butyl ester 9-(2,2,2-trichloro-1,1-dimethylethyl) ester (AK79)*

15

As for bicyclononene AK2, but from bicyclononene AY5 (534 mg, 0.7 mmol), cyclopropyl-[2-(4-methoxyphenoxy)ethyl]amine (228 mg, 1.10 mmol), DIPEA (0.479 mL, 2.8 mmol), DMAP (21 mg, 0.18 mmol), HOBt (113 mg, 0.84 mmol) and EDC·HCl (211 mg, 1.1 mmol) in CH₂Cl₂ (7 mL). Purification by FC yielded the title compound (190 mg, 29%). LC-MS: R_t = 1.26.

(rac.)-(1R, 5S*)-7-{4-[2-(4-Bromophenoxy)ethoxy]phenyl}-6-[cyclopropyl-[2-(3,4-dimethylphenoxy)ethyl]carbamoyl]-3,9-diazabicyclo[3.3.1]non-6-ene-3,9-dicarboxylic acid 3-*tert*-butyl ester 9-(2,2,2-trichloro-1,1-dimethylethyl) ester (AK80)*

As for bicyclononene AK2, but from bicyclononene AY5 (700 mg, 0.918 mmol), cyclopropyl-[2-(3,4-dimethylphenoxy)ethyl]amine (565 mg, 2.75 mmol), DIPEA (0.628 mL, 3.67 mmol), DMAP (28 mg, 0.23 mmol), HOBt (136 mg, 1.01 mmol) and EDC·HCl (264 mg, 1.38 mmol) in CH₂Cl₂ (10 mL). Purification by FC yielded the title compound (199 mg, 23%). LC-MS: R_t = 7.76.

(*rac.*)-(IR*, 5S*)-7-{4-[2-(4-Bromophenoxy)ethoxy]phenyl}-6-{[2-(2-chlorophenyl)ethyl]cyclopropylcarbamoyl}-3,9-diazabicyclo[3.3.1]non-6-ene-3,9-dicarboxylic acid 3-*tert*-butyl ester 9-(2,2,2-trichloro-1,1-dimethylethyl) ester
5 (AK81)

As for bicyclononene AK2, but from bicyclononene AY5 (700 mg, 0.918 mmol), [2-(2-chlorophenyl)ethyl]cyclopropylamine (538 mg, 2.75 mmol), DIPEA (0.628 mL, 3.67 mmol), DMAP (28 mg, 0.23 mmol), HOBr (136 mg, 1.01 mmol) and
10 EDC·HCl (264 mg, 1.38 mmol) in CH₂Cl₂ (10 mL). Purification by FC yielded the title compound (256 mg, 30%). LC-MS: R_t = 7.70.

(*rac.*)-(IR*, 5S*)-7-{4-[2-(4-Bromophenoxy)ethoxy]phenyl}-6-{cyclopropyl-[2-(2,3-difluorophenyl)ethyl]carbamoyl}-3,9-diazabicyclo[3.3.1]non-6-ene-3,9-dicarboxylic acid 3-*tert*-butyl ester 9-(2,2,2-trichloro-1,1-dimethylethyl) ester
15 (AK82)

As for bicyclononene AK2, but from bicyclononene AY5 (700 mg, 0.918 mmol), cyclopropyl-[2-(2,3-difluorophenyl)ethyl]amine (542 mg, 2.75 mmol), DIPEA (0.628 mL, 3.67 mmol), DMAP (28 mg, 0.23 mmol), HOBr (136 mg, 1.01 mmol) and EDC·HCl (264 mg, 1.38 mmol) in CH₂Cl₂ (10 mL). Purification by FC yielded the title compound (245 mg, 28%). LC-MS: R_t = 7.55.

(*rac.*)-(IR*, 5S*)-7-{4-[2-(4-Bromophenoxy)ethoxy]phenyl}-6-{cyclopropyl-[2-(4-fluorophenyl)ethyl]carbamoyl}-3,9-diazabicyclo[3.3.1]non-6-ene-3,9-dicarboxylic acid 3-*tert*-butyl ester 9-(2,2,2-trichloro-1,1-dimethylethyl) ester
25 (AK83)

As for bicyclononene AK2, but from bicyclononene AY5 (700 mg, 0.918 mmol),
30 cyclopropyl-[2-(4-fluorophenyl)ethyl]amine (493 mg, 2.75 mmol), DIPEA (0.628 mL, 3.67 mmol), DMAP (28 mg, 0.23 mmol), HOBr (136 mg, 1.01 mmol) and

EDC·HCl (264 mg, 1.38 mmol) in CH₂Cl₂ (10 mL). Purification by FC yielded the title compound (220 mg, 26%). LC-MS: R_t = 7.51.

5 *(rac.)-(IR*, 5S*)-7-{4-[2-(4-Bromophenoxy)ethoxy]phenyl}-6-[cyclopropyl-(2-o-tolylethyl)carbamoyl]-3,9-diazabicyclo[3.3.1]non-6-ene-3,9-dicarboxylic acid 3-tert-butyl ester 9-(2,2,2-trichloro-1,1-dimethylethyl) ester (AK84)*

As for bicyclononene AK2, but from bicyclononene AY5 (700 mg, 0.918 mmol), cyclopropyl-(2-o-tolylethyl)amine (482 mg, 2.75 mmol), DIPEA (0.628 mL, 3.67 mmol), DMAP (28 mg, 0.23 mmol), HOBr (136 mg, 1.01 mmol) and EDC·HCl (264 mg, 1.38 mmol) in CH₂Cl₂ (10 mL). Purification by FC yielded the title compound (252 mg, 30%). LC-MS: R_t = 7.66.

15 *(rac.)-(IR*, 5S*)-7-{4-[2-(4-Bromophenoxy)ethoxy]phenyl}-6-[cyclopropyl-(3,5-dimethoxybenzyl)carbamoyl]-3,9-diazabicyclo[3.3.1]non-6-ene-3,9-dicarboxylic acid 3-tert-butyl ester 9-(2,2,2-trichloro-1,1-dimethylethyl) ester (AK85)*

20 As for bicyclononene AK2, but from bicyclononene AY5 (700 mg, 0.918 mmol), Cyclopropyl-(3,5-dimethoxy-benzyl)-amine (570 mg, 2.75 mmol), DIPEA (0.628 mL, 3.67 mmol), DMAP (28 mg, 0.23 mmol), HOBr (136 mg, 1.01 mmol) and EDC·HCl (264 mg, 1.38 mmol) in CH₂Cl₂ (10 mL). Purification by FC yielded the title compound (242 mg, 28%). LC-MS: R_t = 7.42.

25 *(rac.)-(IR*, 5S*)-7-{4-[2-(4-Bromophenoxy)ethoxy]phenyl}-6-[cyclopropyl-(2-p-tolylethyl)carbamoyl]-3,9-diazabicyclo[3.3.1]non-6-ene-3,9-dicarboxylic acid 3-tert-butyl ester 9-(2,2,2-trichloro-1,1-dimethylethyl) ester (AK86)*

30 As for bicyclononene AK2, but from bicyclononene AY5 (700 mg, 0.918 mmol), cyclopropyl-(2-p-tolylethyl)amine (482 mg, 2.75 mmol), DIPEA (0.628 mL, 3.67 mmol), DMAP (28 mg, 0.23 mmol), HOBr (136 mg, 1.01 mmol) and EDC·HCl

(264 mg, 1.38 mmol) in CH₂Cl₂ (10 mL). Purification by FC yielded the title compound (246 mg, 29%). LC-MS: R_t = 7.66.

(*rac.*)-(IR*, 5S*)-7-{4-[2-(4-Bromophenoxy)ethoxy]phenyl}-6-{cyclopropyl-[2-5 (2-hydroxyethyl)benzyl]carbamoyl}-3,9-diazabicyclo[3.3.1]non-6-ene-3,9-dicarboxylic acid 3-*tert*-butyl ester 9-(2,2,2-trichloro-1,1-dimethylethyl) ester (AK87)

As for bicyclononene AK2, but from bicyclononene AY5 (1.00 g, 1.31 mmol), (2-10 allylbenzyl)cyclopropylamine (752 mg, 4.02 mmol), DIPEA (0.918 mL, 5.26 mmol), DMAP (41 mg, 0.34 mmol), HOBr (199 mg, 1.47 mmol) and EDC·HCl (385 mg, 2.01 mmol) in CH₂Cl₂ (15 mL). Purification by FC yielded the intermediate compound (875 mg, 72%). R_f = 0.45 (EtOAc/heptane 1:1). LC-MS: R_t = 7.69.

15 Then as for compound AT, but from the former intermediate compound (875 mg, 0.939 mmol), NMO·H₂O (152 mg, 1.13 mmol), and OsO₄ (2.5% in *tert*-BuOH, 0.236 mL, 0.019 mmol) in THF (8 mL), *tert*-BuOH (4 mL) and water (2 mL). Purification of the residue by FC (EtOAc/heptane 1:1 → EtOAc → MeOH/EtOAc 1:9) yielded the 2nd intermediate compound (310 mg, 34%). R_f = 0.05 (EtOAc/heptane 1:1). LC-MS: R_t = 6.86.

Then as for compound AU, but from the 2nd intermediate compound (310 mg, 0.321 mmol) and NaIO₄ (139 mg, 0.481 mmol) in THF (6 mL) and water (2 mL). Drying the residue under high vacuum yielded the 3rd intermediate compound (239 mg, 80%) that was used without further purification. LC-MS: R_t = 7.29.

25 Finally as for compound AV, but from the 3rd intermediate compound (239 mg, 0.256 mmol) and NaBH₄ (11 mg, 0.28 mmol) in MeOH (5 mL). Purification of the residue by FC (EtOAc/heptane 1:4 → 2:3 → 3:2 → 4:1) yielded the title compound (170 mg, 71%). R_f = 0.10 (EtOAc/heptane 1:1). LC-MS: R_t = 7.13.

30 **Compounds of type AL**

(rac.)-(1R, 5S*)-7-{4-[2-(2-Bromo-5-fluorophenoxy)ethyl]phenyl}-6-(methylphenethylcarbamoyl)-3,9-diazabicyclo[3.3.1]non-6-ene-9-carboxylic acid 2,2,2-trichloro-1,1-dimethylethyl ester hydrochloride salt (AL1)*

5 A sol. bicyclononene **AK1** (540 mg, 0.61 mmol) in CH₂Cl₂ (10 mL) was cooled to 0°C. HCl/dioxane (4M, 10 mL) was added and the ice bath was removed. After 4 h stirring at rt the solvents were removed under reduced pressure and the residue dried under high vacuum. The crude was used without further purification.

10

(rac.)-(1R, 5S*)-6-[(2-Chlorobenzyl)cyclopropylcarbamoyl]-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-9-carboxylic acid 2,2,2-trichloro-1,1-dimethyl-ethyl ester hydrochloride salt (AL2)*

15 As for compound **AL1** but from bicyclononene **AK2** (407 mg, 0.47 mmol) in CH₂Cl₂ (5 mL) and HCl/dioxane (4M, 5 mL). LC-MS: R_t = 1.06; ES+: 796.34.

20 *(rac.)-(1R*, 5S*)-6-(Benzylcyclopropylcarbamoyl)-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-9-carboxylic acid 2,2,2-trichloro-1,1-dimethyl-ethyl ester hydrochloride salt (AL3)*

As for compound **AL1** but from bicyclononene **AK3** (570 mg, 0.65 mmol) in CH₂Cl₂ (5 mL) and HCl/dioxane (4M, 5 mL). LC-MS: R_t = 1.04; ES+: 766.34.

25 *(rac.)-(1R*, 5S*)-6-[(2-Chlorobenzyl)ethylcarbamoyl]-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-9-carboxylic acid 2,2,2-trichloro-1,1-dimethylethyl ester hydrochloride salt (AL4)*

30 As for compound **AL1** but from bicyclononene **AK4** (about 0.55 mmol) in CH₂Cl₂ (5 mL) and HCl/dioxane (4M, 5 mL). LC-MS: R_t = 1.19; ES+: 786.25.

(*rac.*)-(1*R*^{*,}, 5*S*^{*})-6-[Cyclopropyl-(2-fluorobenzyl)carbamoyl]-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]-phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-9-carboxylic acid 2,2,2-trichloro-1,1-dimethylethyl ester hydrochloride salt (AL5)

5

As for compound AL1 but from bicyclononene AK5 (about 0.55 mmol) in CH₂Cl₂ (5 mL) and HCl/dioxane (4M, 5 mL). LC-MS: R_t = 1.18; ES+: 782.28.

(*rac.*)-(1*R*^{*,}, 5*S*^{*})-6-[Cyclopropyl-(3-trifluoromethylbenzyl)carbamoyl]-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-9-carboxylic acid 2,2,2-trichloro-1,1-dimethylethyl ester hydrochloride salt (AL6)

As for compound AL1 but from bicyclononene AK6 (about 0.55 mmol) in CH₂Cl₂ (5 mL) and HCl/dioxane (4M, 5 mL). LC-MS: R_t = 1.20.

(*rac.*)-(1*R*^{*,}, 5*S*^{*})-6-[Cyclopropyl-(2-methylbenzyl)carbamoyl]-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-9-carboxylic acid 2,2,2-trichloro-1,1-dimethylethyl ester hydrochloride salt (AL7)

20

As for compound AL1 but from bicyclononene AK7 (about 0.55 mmol) in CH₂Cl₂ (5 mL) and HCl/dioxane (4M, 5 mL). LC-MS: R_t = 1.21; ES+: 778.30.

(*rac.*)-(1*R*^{*,}, 5*S*^{*})-6-[Cyclopropyl-(4-methoxyphenoxy)methyl]carbamoyl]-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-9-carboxylic acid 2,2,2-trichloro-1,1-dimethylethyl ester hydrochloride salt (AL8)

As for compound AL1 but from bicyclononene AK8 (about 0.55 mmol) in CH₂Cl₂ (5 mL) and HCl/dioxane (4M, 5 mL). LC-MS: R_t = 1.21.

(*rac.*)-(IR*, 5S*)-6-[Cyclopropyl-(3-methoxyphenoxy)methyl]carbamoyl]-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-9-carboxylic acid 2,2,2-trichloro-1,1-dimethylethyl ester hydrochloride salt (AL9)

5

As for compound AL1 but from bicyclononene AK9 (about 0.55 mmol) in CH₂Cl₂ (5 mL) and HCl/dioxane (4M, 5 mL). LC-MS: R_t = 1.18.

(*rac.*)-(IR*, 5S*)-6-(Cyclopropyl-*m*-tolyloxymethylcarbamoyl)-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-9-carboxylic acid 2,2,2-trichloro-1,1-dimethylethyl ester hydrochloride salt (AL10)

As for compound AL1 but from bicyclononene AK10 (about 0.55 mmol) in CH₂Cl₂ (5 mL) and HCl/dioxane (4M, 5 mL). LC-MS: R_t = 1.24.

15

(*rac.*)-(IR*, 5S*)-6-[Cyclopropyl-(3,4-dimethylphenoxy)methyl]carbamoyl]-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-9-carboxylic acid 2,2,2-trichloro-1,1-dimethylethyl ester hydrochloride salt (AL11)

20

As for compound AL1 but from bicyclononene AK11 (about 0.55 mmol) in CH₂Cl₂ (5 mL) and HCl/dioxane (4M, 5 mL). LC-MS: R_t = 1.28.

(*rac.*)-(IR*, 5S*)-6-(Cyclopropylphenethylcarbamoyl)-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-9-carboxylic acid 2,2,2-trichloro-1,1-dimethylethyl ester hydrochloride salt (AL12)

As for compound AL1 but from bicyclononene AK12 (about 0.55 mmol) in CH₂Cl₂ (5 mL) and HCl/dioxane (4M, 5 mL). LC-MS: R_t = 1.18; ES+: 778.30.

30

(*rac.*)-(IR*, 5S*)-6-{{[2-(2-Chlorophenyl)ethyl]cyclopropyl}carbamoyl}-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-9-

carboxylic acid 2,2,2-trichloro-1,1-dimethylethyl ester hydrochloride salt
(AL13)

As for compound AL1 but from bicyclononene AK13 (about 0.55 mmol) in
5 CH₂Cl₂ (5 mL) and HCl/dioxane (4M, 5 mL). LC-MS: R_t = 1.21.

(*rac.*)-(1*R*^{*,} 5*S*^{*})-6-{Cyclopropyl-[2-(2,3-difluorophenyl)ethyl]carbamoyl}-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-9-carboxylic acid 2,2,2-trichloro-1,1-dimethylethyl ester (AL14)

10 As for compound AL1 but from bicyclononene AK14 (about 0.55 mmol) in
CH₂Cl₂ (5 mL) and HCl/dioxane (4M, 5 mL). LC-MS: R_t = 1.21.

15 (*rac.*)-(1*R*^{*,} 5*S*^{*})-6-{Cyclopropyl-[2-(4-fluorophenyl)ethyl]carbamoyl}-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-9-carboxylic acid 2,2,2-trichloro-1,1-dimethylethyl ester hydrochloride salt
(AL15)

20 As for compound AL1 but from bicyclononene AK15 (about 0.55 mmol) in
CH₂Cl₂ (5 mL) and HCl/dioxane (4M, 5 mL). LC-MS: R_t = 1.18; ES+: 796.28.

25 (*rac.*)-(1*R*^{*,} 5*S*^{*})-6-[Cyclopropyl-(2-*o*-tolylethyl)carbamoyl]-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diaza-bicyclo[3.3.1]non-6-ene-9-carboxylic acid 2,2,2-trichloro-1,1-dimethylethyl ester hydrochloride salt (AL16)

As for compound AL1 but from bicyclononene AK16 (about 0.55 mmol) in
CH₂Cl₂ (5 mL) and HCl/dioxane (4M, 5 mL). LC-MS: R_t = 1.21; ES+: 794.30.

30 (*rac.*)-(1*R*^{*,} 5*S*^{*})-6-[Cyclopropyl-(3,5-dimethoxybenzyl)carbamoyl]-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-9-carboxylic acid 2,2,2-trichloro-1,1-dimethylethyl ester hydrochloride salt
(AL17)

As for compound **AL1** but from bicyclononene **AK17** (about 0.55 mmol) in CH₂Cl₂ (5 mL) and HCl/dioxane (4M, 5 mL). LC-MS: R_t = 1.16.

- 5 **(rac.)-(1*R**, 5*S**)-6-[Cyclopropyl-(2-*p*-tolylethyl)carbamoyl]-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-9-carboxylic acid 2,2,2-trichloro-1,1-dimethylethyl ester hydrochloride salt (AL18)**

10 As for compound **AL1** but from bicyclononene **AK18** (about 0.55 mmol) in CH₂Cl₂ (5 mL) and HCl/dioxane (4M, 5 mL). LC-MS: R_t = 1.22; ES+: 792.30.

- 15 **(rac.)-(1*R**, 5*S**)-6-[Cyclopropyl-[2-(2-hydroxyethyl)benzyl]carbamoyl]-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-9-carboxylic acid 2,2,2-trichloro-1,1-dimethylethyl ester hydrochloride salt (AL19)**

As for compound **AL1** but from bicyclononene **AK19** (about 0.55 mmol) in CH₂Cl₂ (5 mL) and HCl/dioxane (4M, 5 mL). LC-MS: R_t = 1.01.

- 20 **(rac.)-(1*R**, 5*S**)-7-{4-[3-(2-Bromo-5-fluorophenoxy)propyl]phenyl}-6-[(2-chlorobenzyl)cyclopropylcarbamoyl]-3,9-diazabicyclo[3.3.1]non-6-ene-9-carboxylic acid 2,2,2-trichloro-1,1-dimethylethyl ester hydrochloride salt (AL20)**

- 25 As for compound **AL1** but from bicyclononene **AK20** (about 0.55 mmol) in CH₂Cl₂ (5 mL) and HCl/dioxane (4M, 5 mL). LC-MS: R_t = 5.26; ES+: 806.26.

- 30 **(rac.)-(1*R**, 5*S**)-6-(Benzylcyclopropylcarbamoyl)-7-{4-[3-(2-bromo-5-fluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-9-carboxylic acid 2,2,2-trichloro-1,1-dimethylethyl ester hydrochloride salt (AL21)**

As for compound **AL1** but from bicyclononene **AK21** (519 mg, 0.54 mmol) in CH₂Cl₂ (5 mL) and HCl/dioxane (4M, 5 mL). LC-MS: R_t = 1.06.

5 *(rac.)-(IR*, 5S*)-7-{4-[3-(2-Bromo-5-fluorophenoxy)propyl]phenyl}-6-[(2-chlorobenzyl)ethylcarbamoyl]-3,9-diazabicyclo[3.3.1]non-6-ene-9-carboxylic acid 2,2,2-trichloro-1,1-dimethylethyl ester hydrochloride salt (AL22)*

As for compound **AL1** but from bicyclononene **AK22** (about 0.8 mmol) in CH₂Cl₂ (5 mL) and HCl/dioxane (4M, 5 mL). LC-MS: R_t = 5.30; ES+: 828.33.

10

(rac.)-(IR, 5S*)-7-{4-[3-(2-Bromo-5-fluorophenoxy)propyl]phenyl}-6-[cyclopropyl-(2-fluorobenzyl)carbamoyl]-3,9-diazabicyclo[3.3.1]non-6-ene-9-carboxylic acid 2,2,2-trichloro-1,1-dimethylethyl ester hydrochloride salt (AL23)*

15

As for compound **AL1** but from bicyclononene **AK23** (about 0.8 mmol) in CH₂Cl₂ (5 mL) and HCl/dioxane (4M, 5 mL).

20

(rac.)-(IR, 5S*)-7-{4-[3-(2-Bromo-5-fluorophenoxy)propyl]phenyl}-6-[cyclopropyl-(3-trifluoromethylbenzyl)carbamoyl]-3,9-diazabicyclo[3.3.1]non-6-ene-9-carboxylic acid 2,2,2-trichloro-1,1-dimethylethyl ester hydrochloride salt (AL24)*

As for compound **AL1** but from bicyclononene **AK24** (about 0.8 mmol) in CH₂Cl₂ (5 mL) and HCl/dioxane (4M, 5 mL). LC-MS: R_t = 5.33; ES+: 820.40.

25

(rac.)-(IR, 5S*)-7-{4-[3-(2-Bromo-5-fluorophenoxy)propyl]phenyl}-6-[cyclopropyl-(2-methylbenzyl)carbamoyl]-3,9-diazabicyclo[3.3.1]non-6-ene-9-carboxylic acid 2,2,2-trichloro-1,1-dimethylethyl ester hydrochloride salt (AL25)*

30

As for compound **AL1** but from bicyclononene **AK25** (about 0.8 mmol) in CH₂Cl₂ (5 mL) and HCl/dioxane (4M, 5 mL). LC-MS: R_t = 1.06.

(*rac.*)-(IR*, 5S*)-7-{4-[3-(2-Bromo-5-fluorophenoxy)propyl]phenyl}-6-{cyclo-
propyl-[2-(4-methoxyphenoxy)ethyl]carbamoyl}-3,9-diazabicyclo[3.3.1]non-
6-ene-9-carboxylic acid 2,2,2-trichloro-1,1-dimethylethyl ester hydrochloride
5 salt (AL26)

As for compound AL1 but from bicyclononene AK26 (about 0.8 mmol) in CH₂Cl₂ (5 mL) and HCl/dioxane (4M, 5 mL). LC-MS: R_t = 5.08; ES+: 866.40.

10 (*rac.*)-(IR*, 5S*)-7-{4-[3-(2-Bromo-5-fluorophenoxy)propyl]phenyl}-6-[cyclo-
propyl-(2-*m*-tolyloxyethyl)carbamoyl]-3,9-diazabicyclo[3.3.1]non-6-ene-9-
carboxylic acid 2,2,2-trichloro-1,1-dimethylethyl ester hydrochloride salt
(AL27)

15 As for compound AL1 but from bicyclononene AK27 (about 0.8 mmol) in CH₂Cl₂ (5 mL) and HCl/dioxane (4M, 5 mL). LC-MS: R_t = 1.07.

20 (*rac.*)-(IR*, 5S*)-7-{4-[3-(2-Bromo-5-fluorophenoxy)propyl]phenyl}-6-{cyclo-
propyl-[2-(3,4-dimethylphenoxy)ethyl]carbamoyl}-3,9-diazabicyclo[3.3.1]-
non-6-ene-9-carboxylic acid 2,2,2-trichloro-1,1-dimethylethyl ester hydrochloride salt (AL28)

25 As for compound AL1 but from bicyclononene AK28 (about 0.8 mmol) in CH₂Cl₂ (5 mL) and HCl/dioxane (4M, 5 mL). LC-MS: R_t = 1.08.

-(IR*, 5S*)-7-{4-[3-(2-Bromo-5-fluorophenoxy)propyl]phenyl}-6-(cyclo-
propylphenethylcarbamoyl)-3,9-diazabicyclo[3.3.1]non-6-ene-9-carboxylic
acid 2,2,2-trichloro-1,1-dimethylethyl ester hydrochloride salt (AL29)

30 As for compound AL1 but from bicyclononene AK29 (about 0.8 mmol) in CH₂Cl₂ (5 mL) and HCl/dioxane (4M, 5 mL). LC-MS: R_t = 5.25; ES+: 820.38.

(*rac.*)-(IR*, 5S*)-7-{4-[3-(2-Bromo-5-fluorophenoxy)propyl]phenyl}-6-{[2-(2-chlorophenyl)ethyl]cyclopropylcarbamoyl}-3,9-diazabicyclo[3.3.1]non-6-ene-9-carboxylic acid 2,2,2-trichloro-1,1-dimethylethyl ester hydrochloride salt (AL30)

5

As for compound AL1 but from bicyclononene AK30 (about 0.8 mmol) in CH₂Cl₂ (5 mL) and HCl/dioxane (4M, 5 mL). LC-MS: R_t = 5.35; ES+: 854.30.

(*rac.*)-(IR*, 5S*)-7-{4-[3-(2-Bromo-5-fluorophenoxy)propyl]phenyl}-6-{cyclo-10 propyl-[2-(2,3-difluorophenyl)ethyl]carbamoyl}-3,9-diazabicyclo[3.3.1]non-6-ene-9-carboxylic acid 2,2,2-trichloro-1,1-dimethylethyl ester hydrochloride salt (AL31)

As for compound AL1 but from bicyclononene AK31 (about 0.8 mmol) in 15 CH₂Cl₂ (5 mL) and HCl/dioxane (4M, 5 mL). LC-MS: R_t = 5.40; ES+: 856.38.

(*rac.*)-(IR*, 5S*)-7-{4-[3-(2-Bromo-5-fluorophenoxy)propyl]phenyl}-6-{cyclo-15 propyl-[2-(4-fluorophenyl)ethyl]carbamoyl}-3,9-diazabicyclo[3.3.1]non-6-ene-9-carboxylic acid 2,2,2-trichloro-1,1-dimethylethyl ester hydrochloride 20 salt (AL32)

As for compound AL1 but from bicyclononene AK32 (about 0.8 mmol) in CH₂Cl₂ (5 mL) and HCl/dioxane (4M, 5 mL). LC-MS: R_t = 5.28; ES+: 838.40.

25 (*rac.*)-(IR*, 5S*)-7-{4-[3-(2-Bromo-5-fluorophenoxy)propyl]phenyl}-6-[cyclo-propyl-(2-*o*-tolylethyl)carbamoyl]-3,9-diazabicyclo[3.3.1]non-6-ene-9-carbo-30 xylic acid 2,2,2-trichloro-1,1-dimethylethyl ester hydrochloride salt (AL33)

As for compound AL1 but from bicyclononene AK33 (about 0.8 mmol) in 30 CH₂Cl₂ (5 mL) and HCl/dioxane (4M, 5 mL). LC-MS: R_t = 5.36; ES+: 834.42.

1:1-Mixture of (*rac.*)-(1*R*^{*, 5*S*})-7-{4-[3-(2-Bromo-5-fluorophenoxy)propyl]-phenyl}-6-[((2*R*^{*})-2-hydroxy-2-phenylethyl)methylcarbamoyl]-3,9-diaza-bicyclo[3.3.1]non-6-ene-9-carboxylic acid 2,2,2-trichloro-1,1-dimethylethyl ester hydrochloride salt and (*rac.*)-(1*R*^{*, 5*S*})-7-{4-[3-(2-Bromo-5-fluorophenoxy)propyl]phenyl}-6-[((2*S*^{*})-2-hydroxy-2-phenylethyl)methylcarbamoyl]-3,9-diazabicyclo[3.3.1]non-6-ene-9-carboxylic acid 2,2,2-trichloro-1,1-dimethylethyl ester hydrochloride salt (AL34)

As for compound AL1 but from bicyclononenes AK34 (about 0.3 mmol) in
10 CH₂Cl₂ (5 mL) and HCl/dioxane (4M, 5 mL). LC-MS: R_t = 0.99.

(*rac.*)-(1*R*^{*, 5*S*})-7-{4-[3-(2-Bromo-5-fluorophenoxy)propyl]phenyl}-6-[cyclo-propyl-(3,5-dimethoxybenzyl)carbamoyl]-3,9-diazabicyclo[3.3.1]non-6-ene-9-carboxylic acid 2,2,2-trichloro-1,1-dimethylethyl ester hydrochloride salt
15 (AL35)

As for compound AL1 but from bicyclononene AK35 (about 0.8 mmol) in
CH₂Cl₂ (5 mL) and HCl/dioxane (4M, 5 mL). LC-MS: R_t = 5.23; ES+: 866.40.

20 (*rac.*)-(1*R*^{*, 5*S*})-7-{4-[3-(2-Bromo-5-fluorophenoxy)propyl]phenyl}-6-[cyclo-propyl-(2-*p*-tolylethyl)carbamoyl]-3,9-diazabicyclo[3.3.1]non-6-ene-9-carboxylic acid 2,2,2-trichloro-1,1-dimethylethyl ester hydrochloride salt (AL36)

As for compound AL1 but from bicyclononene AK36 (about 0.8 mmol) in
25 CH₂Cl₂ (5 mL) and HCl/dioxane (4M, 5 mL). LC-MS: R_t = 5.59; ES+: 834.42.

(*rac.*)-(1*R*^{*, 5*S*})-7-{4-[3-(2-Bromo-5-fluorophenoxy)propyl]phenyl}-6-{cyclo-propyl-[2-(2-hydroxyethyl)benzyl]carbamoyl}-3,9-diazabicyclo[3.3.1]non-6-ene-9-carboxylic acid 2,2,2-trichloro-1,1-dimethylethyl ester hydrochloride
30 salt (AL37)

As for compound **AL1** but from bicyclononene **AK37** (about 0.8 mmol) in CH₂Cl₂ (5 mL) and HCl/dioxane (4M, 5 mL). LC-MS: R_t = 1.01.

5 *(rac.)-(1R*, 5S*)-6-[(2-Chlorobenzyl)cyclopropylcarbamoyl]-7-{4-[2-(2,3,6-trimethylphenoxy)ethyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-9-carboxylic acid 2,2,2-trichloro-1,1-dimethylethyl ester hydrochloride salt (AL38)*

As for compound **AL1** but from bicyclononene **AK38** (312 mg, 0.35 mmol) in CH₂Cl₂ (5 mL) and HCl/dioxane (4M, 5 mL). LC-MS: R_t = 1.08; ES+: 774.33.

10

(rac.)-(1R, 5S*)-6-(Benzylcyclopropylcarbamoyl)-7-{4-[2-(2,3,6-trimethylphenoxy)ethyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-9-carboxylic acid 2,2,2-trichloro-1,1-dimethylethyl ester hydrochloride salt (AL39)*

15 As for compound **AL1** but from bicyclononene **AK39** (340 mg, 0.40 mmol) in CH₂Cl₂ (5 mL) and HCl/dioxane (4M, 5 mL). LC-MS: R_t = 1.05; ES+: 740.42.

20 *(rac.)-(1R*, 5S*)-6-[(2-Chlorobenzyl)ethylcarbamoyl]-7-{4-[2-(2,3,6-trimethyl-phenoxy)ethyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-9-carboxylic acid 2,2,2-trichloro-1,1-dimethylethyl ester hydrochloride salt (AL40)*

As for compound **AL1** but from bicyclononene **AK40** (374 mg, 0.43 mmol) in CH₂Cl₂ (5 mL) and HCl/dioxane (4M, 5 mL). LC-MS: R_t = 1.07; ES+: 762.34.

25

(rac.)-(1R, 5S*)-6-[Cyclopropyl-(2-fluorobenzyl)carbamoyl]-7-{4-[2-(2,3,6-trimethylphenoxy)ethyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-9-carboxylic acid 2,2,2-trichloro-1,1-dimethylethyl ester hydrochloride salt (AL41)*

30 As for compound **AL1** but from bicyclononene **AK41** (350 mg, 0.41 mmol) in CH₂Cl₂ (5 mL) and HCl/dioxane (4M, 5 mL). LC-MS: R_t = 1.06; ES+: 758.38.

(*rac.*)-(1*R*^{*,} 5*S*^{*})-6-[Cyclopropyl-(3-trifluoromethylbenzyl)carbamoyl]-7-{4-[2-(2,3,6-trimethylphenoxy)ethyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-9-carboxylic acid 2,2,2-trichloro-1,1-dimethylethyl ester hydrochloride salt (AL42)

5

As for compound AL1 but from bicyclononene AK42 (294 mg, 0.32 mmol) in CH₂Cl₂ (5 mL) and HCl/dioxane (4M, 5 mL). LC-MS: R_t = 1.08.

(*rac.*)-(1*R*^{*,} 5*S*^{*})-6-[Cyclopropyl-(2-methylbenzyl)carbamoyl]-7-{4-[2-(2,3,6-trimethylphenoxy)ethyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-9-carboxylic acid 2,2,2-trichloro-1,1-dimethylethyl ester hydrochloride salt (AL43)

10 As for compound AL1 but from bicyclononene AK43 (322 mg, 0.38 mmol) in CH₂Cl₂ (5 mL) and HCl/dioxane (4M, 5 mL). LC-MS: R_t = 1.08; ES+: 752.39.

15

(*rac.*)-(1*R*^{*,} 5*S*^{*})-6-[Cyclopropyl-[2-(4-methoxyphenoxy)ethyl]carbamoyl]-7-{4-[2-(2,3,6-trimethylphenoxy)ethyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-9-carboxylic acid 2,2,2-trichloro-1,1-dimethylethyl ester hydrochloride salt (AL44)

20

As for compound AL1 but from bicyclononene AK44 (159 mg, 0.18 mmol) in CH₂Cl₂ (5 mL) and HCl/dioxane (4M, 5 mL). LC-MS: R_t = 1.07.

(*rac.*)-(1*R*^{*,} 5*S*^{*})-6-[Cyclopropyl-[2-(3-methoxyphenoxy)ethyl]carbamoyl]-7-{4-[2-(2,3,6-trimethylphenoxy)ethyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-9-carboxylic acid 2,2,2-trichloro-1,1-dimethylethyl ester (AL45)

25 As for compound AL1 but from bicyclononene AK45 (237 mg, 0.26 mmol) in CH₂Cl₂ (5 mL) and HCl/dioxane (4M, 5 mL). LC-MS: R_t = 1.00.

30

(*rac.*)-(1*R*^{*,} 5*S*^{*})-6-[Cyclopropyl-(2-*m*-tolyloxyethyl)carbamoyl]-7-{4-[2-(2,3,6-trimethylphenoxy)ethyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-9-

**carboxylic acid 2,2,2-trichloro-1,1-dimethylethyl ester hydrochloride salt
(AL46)**

As for compound AL1 but from bicyclononene AK46 (185 mg, 0.21 mmol) in
5 CH₂Cl₂ (5 mL) and HCl/dioxane (4M, 5 mL). LC-MS: R_t = 1.09; ES+: 784.40.

**(rac.)-(IR*, 5S*)-6-(Cyclopropylphenethylcarbamoyl)-7-{4-[2-(2,3,6-trimethyl-phenoxy)ethyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-9-carboxylic acid 2,2,2-trichloro-1,1-dimethylethyl ester hydrochloride salt
(AL47)**

As for compound AL1 but from bicyclononene AK47 (about 0.3 mmol) in
CH₂Cl₂ (5 mL) and HCl/dioxane (4M, 5 mL). LC-MS: R_t = 1.09; ES+: 754.44.

**15 (rac.)-(IR*, 5S*)-6-{[2-(2-Chlorophenyl)ethyl]cyclopropylcarbamoyl}-7-{4-[2-(2,3,6-trimethylphenoxy)ethyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-9-carboxylic acid 2,2,2-trichloro-1,1-dimethylethyl ester hydrochloride salt
(AL48)**

20 As for compound AL1 but from bicyclononene AK48 (about 0.3 mmol) in
CH₂Cl₂ (5 mL) and HCl/dioxane (4M, 5 mL). LC-MS: R_t = 1.10; ES+: 788.41.

**25 (rac.)-(IR*, 5S*)-6-{Cyclopropyl-[2-(2,3-difluorophenyl)ethyl]carbamoyl}-7-{4-[2-(2,3,6-trimethylphenoxy)ethyl]-phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-9-carboxylic acid 2,2,2-trichloro-1,1-dimethylethyl ester hydrochloride salt
(AL49)**

As for compound AL1 but from bicyclononene AK49 (about 0.3 mmol) in
CH₂Cl₂ (5 mL) and HCl/dioxane (4M, 5 mL). LC-MS: R_t = 1.09; ES+: 788.41.

30

(rac.)-(IR*, 5S*)-6-{Cyclopropyl-[2-(4-fluorophenyl)ethyl]carbamoyl}-7-{4-[2-(2,3,6-trimethylphenoxy)ethyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-9-

**carboxylic acid 2,2,2-trichloro-1,1-dimethylethyl ester hydrochloride salt
(AL50)**

As for compound **AL1** but from bicyclononene **AK50** (about 0.3 mmol) in
5 CH₂Cl₂ (5 mL) and HCl/dioxane (4M, 5 mL). LC-MS: R_t = 1.09; ES+: 772.41.

(rac.)-(1*R*^{*}, 5*S*^{*})-6-[Cyclopropyl-(2-*o*-tolylethyl)carbamoyl]-7-{4-[2-(2,3,6-trimethylphenoxy)ethyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-9-carboxylic acid 2,2,2-trichloro-1,1-dimethylethyl ester hydrochloride salt (AL51)

10

As for compound **AL1** but from bicyclononene **AK51** (about 0.3 mmol) in CH₂Cl₂ (5 mL) and HCl/dioxane (4M, 5 mL). LC-MS: R_t = 1.10; ES+: 768.44.

15 **1:1-Mixture of (rac.)-(1*R*^{*}, 5*S*^{*})-6-[(2*S*^{*})-2-hydroxy-2-phenylethyl)methylcarbamoyl]-7-{4-[2-(2,3,6-trimethylphenoxy)ethyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-9-carboxylic acid 2,2,2-trichloro-1,1-dimethylethyl ester hydrochloride salt and (rac.)-(1*R*^{*}, 5*S*^{*})-6-[(2*S*^{*})-2-hydroxy-2-phenylethyl)methylcarbamoyl]-7-{4-[2-(2,3,6-trimethylphenoxy)ethyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-9-carboxylic acid 2,2,2-trichloro-1,1-dimethylethyl ester hydrochloride salt (AL52)**

20

As for compound **AL1** but from bicyclononene **AK52** (about 0.3 mmol) in CH₂Cl₂ (5 mL) and HCl/dioxane (4M, 5 mL). LC-MS: R_t = 1.02; ES+: 744.44.

25 **(rac.)-(1*R*^{*}, 5*S*^{*})-6-[Cyclopropyl-(3,5-dimethoxybenzyl)carbamoyl]-7-{4-[2-(2,3,6-trimethylphenoxy)ethyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-9-carboxylic acid 2,2,2-trichloro-1,1-dimethylethyl ester hydrochloride salt (AL53)**

30 As for compound **AL1** but from bicyclononene **AK53** (about 0.3 mmol) in CH₂Cl₂ (5 mL) and HCl/dioxane (4M, 5 mL). LC-MS: R_t = 1.07.

(rac.)-(IR, 5S*)-6-[Cyclopropyl-(2-p-tolylethyl)carbamoyl]-7-{4-[2-(2,3,6-trimethylphenoxy)ethyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-9-carboxylic acid 2,2,2-trichloro-1,1-dimethylethyl ester hydrochloride salt (AL54)*

5 As for compound AL1 but from bicyclononene AK54 (about 0.3 mmol) in CH₂Cl₂ (5 mL) and HCl/dioxane (4M, 5 mL). LC-MS: R_t = 1.10; ES+: 768.44.

10 *(rac.)-(IR*, 5S*)-6-[Cyclopropyl-[2-(2-hydroxyethyl)benzyl]carbamoyl]-7-{4-[2-(2,3,6-trimethylphenoxy)ethyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-9-carboxylic acid 2,2,2-trichloro-1,1-dimethylethyl ester hydrochloride salt (AL55)*

15 As for compound AL1 but from bicyclononene AK55 (about 0.3 mmol) in CH₂Cl₂ (5 mL) and HCl/dioxane (4M, 5 mL). LC-MS: R_t = 1.03; ES+: 784.44.

20 *(rac.)-(IR*, 5S*)-6-[(2-Chlorobenzyl)cyclopropylcarbamoyl]-7-{4-[2-(2-chloro-4,5-dimethylphenoxy)ethoxy]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-9-carboxylic acid 2,2,2-trichloro-1,1-dimethylethyl ester hydrochloride salt (AL56)*

25 As for compound AL1 but from bicyclononene AK56 (about 0.5 mmol) in CH₂Cl₂ (5 mL) and HCl/dioxane (4M, 5 mL). LC-MS: R_t = 1.08; ES+: 785.26.

30 *(rac.)-(IR*, 5S*)-6-[(2-Chlorobenzyl)ethylcarbamoyl]-7-{4-[2-(2-chloro-4,5-dimethylphenoxy)ethoxy]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-9-carboxylic acid 2,2,2-trichloro-1,1-dimethylethyl ester hydrochloride salt (AL57)*

As for compound AL1 but from bicyclononene AK57 (about 0.5 mmol) in CH₂Cl₂ (5 mL) and HCl/dioxane (4M, 5 mL). LC-MS: R_t = 1.06.

(rac.)-(IR, 5S*)-7-{4-[2-(2-Chloro-4,5-dimethylphenoxy)ethoxy]phenyl}-6-[cyclopropyl-(2-fluorobenzyl)carbamoyl]-3,9-diazabicyclo[3.3.1]non-6-ene-9-*

carboxylic acid 2,2,2-trichloro-1,1-dimethylethyl ester hydrochloride salt
(AL58)

As for compound AL1 but from bicyclononene AK58 (about 0.5 mmol) in
5 CH₂Cl₂ (5 mL) and HCl/dioxane (4M, 5 mL). LC-MS: R_t = 1.06.

(*rac.*)-(IR*, 5S*)-7-{4-[2-(2-Chloro-4,5-dimethylphenoxy)ethoxy]phenyl}-6-[cyclopropyl-(3-trifluoromethylbenzyl)carbamoyl]-3,9-diazabicyclo[3.3.1]-non-6-ene-9-carboxylic acid 2,2,2-trichloro-1,1-dimethylethyl ester
10 hydrochloride salt (AL59)

As for compound AL1 but from bicyclononene AK59 (about 0.5 mmol) in
CH₂Cl₂ (5 mL) and HCl/dioxane (4M, 5 mL). LC-MS: R_t = 1.08.

15 (*rac.*)-(IR*, 5S*)-7-{4-[2-(2-Chloro-4,5-dimethylphenoxy)ethoxy]phenyl}-6-[cyclopropyl-(2-methylbenzyl)carbamoyl]-3,9-diazabicyclo[3.3.1]non-6-ene-9-carboxylic acid 2,2,2-trichloro-1,1-dimethylethyl ester hydrochloride salt
(AL60)

20 As for compound AL1 but from bicyclononene AK60 (about 0.5 mmol) in
CH₂Cl₂ (5 mL) and HCl/dioxane (4M, 5 mL). LC-MS: R_t = 1.07; ES+: 788.40.

(*rac.*)-(IR*, 5S*)-7-{4-[2-(2-Chloro-4,5-dimethylphenoxy)ethoxy]phenyl}-6-[cyclopropyl-[2-(4-methoxyphenoxy)ethyl]carbamoyl]-3,9-diazabicyclo[3.3.1]non-6-ene-9-carboxylic acid 2,2,2-trichloro-1,1-dimethylethyl ester
25 hydrochloride salt (AL61)

As for compound AL1 but from bicyclononene AK61 (about 0.5 mmol) in
CH₂Cl₂ (5 mL) and HCl/dioxane (4M, 5 mL). LC-MS: R_t = 1.06.

30

(*rac.*)-(IR*, 5S*)-7-{4-[2-(2-Chloro-4,5-dimethylphenoxy)ethoxy]phenyl}-6-[cyclopropyl-[2-(3-methoxyphenoxy)ethyl]carbamoyl]-3,9-

diazabicyclo[3.3.1]non-6-ene-9-carboxylic acid 2,2,2-trichloro-1,1-dimethylethyl ester hydrochloride salt (AL62)

As for compound **AL1** but from bicyclononene **AK62** (about 0.5 mmol) in
5 CH₂Cl₂ (5 mL) and HCl/dioxane (4M, 5 mL). LC-MS: R_t = 1.07.

(rac.)-(IR*, 5S*)-7-{4-[2-(2-Chloro-4,5-dimethylphenoxy)ethoxy]phenyl}-6-[cyclopropyl-(2-m-tolyloxyethyl)carbamoyl]-3,9-diazabicyclo[3.3.1]non-6-ene-9-carboxylic acid 2,2,2-trichloro-1,1-dimethylethyl ester hydrochloride salt (AL63)

As for compound **AL1** but from bicyclononene **AK63** (about 0.5 mmol) in
CH₂Cl₂ (5 mL) and HCl/dioxane (4M, 5 mL). LC-MS: R_t = 1.08.

15 (rac.)-(IR*, 5S*)-7-{4-[2-(2-Chloro-4,5-dimethylphenoxy)ethoxy]phenyl}-6-[cyclopropylphenethylcarbamoyl]-3,9-diazabicyclo[3.3.1]non-6-ene-9-carboxylic acid 2,2,2-trichloro-1,1-dimethylethyl ester hydrochloride salt (AL64)

As for compound **AL1** but from bicyclononene **AK64** (about 0.5 mmol) in
20 CH₂Cl₂ (5 mL) and HCl/dioxane (4M, 5 mL). LC-MS: R_t = 1.07; ES+: 788.39.

(rac.)-(IR*, 5S*)-7-{4-[2-(2-Chloro-4,5-dimethylphenoxy)ethoxy]phenyl}-6-[[2-(2-chlorophenyl)ethyl]cyclopropylcarbamoyl]-3,9-diazabicyclo[3.3.1]non-6-ene-9-carboxylic acid 2,2,2-trichloro-1,1-dimethylethyl ester hydrochloride salt (AL65)

As for compound **AL1** but from bicyclononene **AK65** (about 0.5 mmol) in
CH₂Cl₂ (5 mL) and HCl/dioxane (4M, 5 mL). LC-MS: R_t = 1.08.

30 (rac.)-(IR*, 5S*)-7-{4-[2-(2-Chloro-4,5-dimethylphenoxy)ethoxy]phenyl}-6-[cyclopropyl-[2-(2,3-difluorophenyl)ethyl]carbamoyl]-3,9-diazabicyclo-

[3.3.1]non-6-ene-9-carboxylic acid 2,2,2-trichloro-1,1-dimethylethyl ester hydrochloride salt (AL66)

As for compound AL1 but from bicyclononene AK66 (about 0.5 mmol) in
5 CH₂Cl₂ (5 mL) and HCl/dioxane (4M, 5 mL). LC-MS: R_t = 1.06.

(rac.)-(1*R*^{*}, 5*S*^{*})-7-{4-[2-(2-Chloro-4,5-dimethylphenoxy)ethoxy]phenyl}-6-{cyclopropyl-[2-(4-fluorophenyl)ethyl]carbamoyl}-3,9-diazabicyclo[3.3.1]non-6-ene-9-carboxylic acid 2,2,2-trichloro-1,1-dimethylethyl ester hydrochloride salt (AL67)

As for compound AL1 but from bicyclononene AK67 (about 0.5 mmol) in CH₂Cl₂ (5 mL) and HCl/dioxane (4M, 5 mL). LC-MS: R_t = 1.06.

15 (rac.)-(1*R*^{*}, 5*S*^{*})-7-{4-[2-(2-Chloro-4,5-dimethylphenoxy)ethoxy]phenyl}-6-[cyclopropyl-(2-*o*-tolylethyl)carbamoyl]-3,9-diazabicyclo[3.3.1]non-6-ene-9-carboxylic acid 2,2,2-trichloro-1,1-dimethylethyl ester (AL68)

As for compound AL1 but from bicyclononene AK68 (about 0.5 mmol) in
20 CH₂Cl₂ (5 mL) and HCl/dioxane (4M, 5 mL). LC-MS: R_t = 1.08.

25 1:1-Mixture of (rac.)-(1*R*^{*}, 5*S*^{*})-7-{4-[2-(2-chloro-4,5-dimethylphenoxy)-ethoxy]phenyl}-6-[(2*R*^{*})2-hydroxy-2-phenylethyl)methylcarbamoyl]-3,9-diazabicyclo[3.3.1]non-6-ene-9-carboxylic acid 2,2,2-trichloro-1,1-dimethyl-ethyl ester hydrochloride salt and (rac.)-(1*R*^{*}, 5*S*^{*})-7-{4-[2-(2-chloro-4,5-dimethylphenoxy)ethoxy]phenyl}-6-[(2*S*^{*})2-hydroxy-2-phenylethyl)methylcarbamoyl]-3,9-diazabicyclo[3.3.1]non-6-ene-9-carboxylic acid 2,2,2-trichloro-1,1-dimethylethyl ester hydrochloride salt (AL69)

30 As for compound AL1 but from bicyclononene AK69 (about 0.5 mmol) in CH₂Cl₂ (5 mL) and HCl/dioxane (4M, 5 mL). LC-MS: R_t = 0.99; ES+: 780.37.

(*rac.*)-(1*R*^{*,}, 5*S*^{*})-7-{4-[2-(2-Chloro-4,5-dimethylphenoxy)ethoxy]phenyl}-6-[cyclopropyl-(3,5-dimethoxybenzyl)carbamoyl]-3,9-diazabicyclo[3.3.1]non-6-ene-9-carboxylic acid 2,2,2-trichloro-1,1-dimethylethyl ester hydrochloride salt (AL70)

5

As for compound AL1 but from bicyclononene AK70 (about 0.5 mmol) in CH₂Cl₂ (5 mL) and HCl/dioxane (4M, 5 mL). LC-MS: R_t = 1.06.

(*rac.*)-(1*R*^{*,}, 5*S*^{*})-7-{4-[2-(2-Chloro-4,5-dimethylphenoxy)ethoxy]phenyl}-6-[cyclopropyl-(2-*p*-tolylethyl)carbamoyl]-3,9-diazabicyclo[3.3.1]non-6-ene-9-carboxylic acid 2,2,2-trichloro-1,1-dimethylethyl ester hydrochloride salt (AL71)

As for compound AL1 but from bicyclononene AK71 (about 0.5 mmol) in CH₂Cl₂ (5 mL) and HCl/dioxane (4M, 5 mL). LC-MS: R_t = 1.08.

(*rac.*)-(1*R*^{*,}, 5*S*^{*})-7-{4-[2-(2-Chloro-4,5-dimethylphenoxy)ethoxy]phenyl}-6-[cyclopropyl-[2-(2-hydroxyethyl)benzyl]carbamoyl]-3,9-diazabicyclo[3.3.1]-non-6-ene-9-carboxylic acid 2,2,2-trichloro-1,1-dimethylethyl ester hydrochloride salt

20

As for compound AL1 but from bicyclononene AK72 (about 0.5 mmol) in CH₂Cl₂ (5 mL) and HCl/dioxane (4M, 5 mL). LC-MS: R_t = 1.01.

25 (*rac.*)-(1*R*^{*,}, 5*S*^{*})-7-{4-[2-(4-Bromophenoxy)ethoxy]phenyl}-6-[2-chlorobenzyl)cyclopropylcarbamoyl]-3,9-diazabicyclo[3.3.1]non-6-ene-9-carboxylic acid 2,2,2-trichloro-1,1-dimethyl-ethyl ester hydrochloride salt (AL73)

As for compound AL1 but from bicyclononene AK73 (0.28 mmol) in CH₂Cl₂ (5 mL) and HCl/dioxane (4M, 5 mL). LC-MS: R_t = 4.98; ES+: 824.32.

(rac.)-(1R, 5S*)-6-(Benzylcyclopropylcarbamoyl)-7-{4-[2-(4-bromophenoxy)ethoxy]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-9-carboxylic acid 2,2,2-trichloro-1,1-dimethylethyl ester hydrochloride salt (AL74)*

5 As for compound **AL1** but from bicyclononene **AK74** (0.27 mmol) in CH₂Cl₂ (5 mL) and HCl/dioxane (4M, 5 mL). LC-MS: R_t = 1.03; ES+: 792.36.

10 *(rac.)-(1R*, 5S*)-7-{4-[2-(4-Bromophenoxy)ethoxy]phenyl}-6-[(2-chlorobenzyl)ethylcarbamoyl]-3,9-diazabicyclo[3.3.1]non-6-ene-9-carboxylic acid 2,2,2-trichloro-1,1-dimethylethyl ester hydrochloride salt (AL75)*

As for compound **AL1** but from bicyclononene **AK75** (0.22 mmol) in CH₂Cl₂ (5 mL) and HCl/dioxane (4M, 5 mL). LC-MS: R_t = 4.92; ES+: 812.35.

15 *(rac.)-(1R*, 5S*)-7-{4-[2-(4-Bromophenoxy)ethoxy]phenyl}-6-[cyclopropyl-(2-fluorobenzyl)carbamoyl]-3,9-diazabicyclo[3.3.1]non-6-ene-9-carboxylic acid 2,2,2-trichloro-1,1-dimethylethyl ester hydrochloride salt (AL76)*

20 As for compound **AL1** but from bicyclononene **AK76** (0.26 mmol) in CH₂Cl₂ (5 mL) and HCl/dioxane (4M, 5 mL). LC-MS: R_t = 4.78; ES+: 808.40.

25 *(rac.)-(1R*, 5S*)-7-{4-[2-(4-Bromophenoxy)ethoxy]phenyl}-6-[cyclopropyl-(3-trifluoromethylbenzyl)carbamoyl]-3,9-diazabicyclo[3.3.1]non-6-ene-9-carboxylic acid 2,2,2-trichloro-1,1-dimethylethyl ester hydrochloride salt (AL77)*

As for compound **AL1** but from bicyclononene **AK77** (0.29 mmol) in CH₂Cl₂ (5 mL) and HCl/dioxane (4M, 5 mL). LC-MS: R_t = 4.97; ES+: 858.42.

30 *(rac.)-(1R*, 5S*)-7-{4-[2-(4-Bromophenoxy)ethoxy]phenyl}-6-[cyclopropyl-(2-methylbenzyl)carbamoyl]-3,9-diazabicyclo[3.3.1]non-6-ene-9-carboxylic acid 2,2,2-trichloro-1,1-dimethylethyl ester hydrochloride salt (AL78)*

As for compound **AL1** but from bicyclononene **AK78** (0.28 mmol) in CH₂Cl₂ (5 mL) and HCl/dioxane (4M, 5 mL). LC-MS: R_t = 5.01; ES+: 804.42.

5 *(rac.)-(IR*, 5S*)-7-{4-[2-(4-Bromophenoxy)ethoxy]phenyl}-6-{cyclopropyl-[2-(4-methoxyphenoxy)ethyl]carbamoyl}-3,9-diazabicyclo[3.3.1]non-6-ene-9-carboxylic acid 2,2,2-trichloro-1,1-dimethylethyl ester hydrochloride salt (AL79)*

10 As for compound **AL1** but from bicyclononene **AK79** (0.2 mmol) in CH₂Cl₂ (5 mL) and HCl/dioxane (4M, 5 mL). LC-MS: R_t = 5.03; ES+: 850.44.

15 *(rac.)-(IR*, 5S*)-7-{4-[2-(4-Bromophenoxy)ethoxy]phenyl}-6-{cyclopropyl-[2-(3,4-dimethylphenoxy)ethyl]carbamoyl}-3,9-diazabicyclo[3.3.1]non-6-ene-9-carboxylic acid 2,2,2-trichloro-1,1-dimethylethyl ester hydrochloride salt (AL80)*

As for compound **AL1** but from bicyclononene **AK80** (0.21 mmol) in CH₂Cl₂ (5 mL) and HCl/dioxane (4M, 5 mL). LC-MS: R_t = 5.11; ES+: 848.43.

20 *(rac.)-(IR*, 5S*)-7-{4-[2-(4-Bromophenoxy)ethoxy]phenyl}-6-{[2-(2-chlorophenyl)ethyl]cyclopropylcarbamoyl}-3,9-diazabicyclo[3.3.1]non-6-ene-9-carboxylic acid 2,2,2-trichloro-1,1-dimethylethyl ester hydrochloride salt (AL81)*

25 As for compound **AL1** but from bicyclononene **AK81** (0.27 mmol) in CH₂Cl₂ (5 mL) and HCl/dioxane (4M, 5 mL). LC-MS: R_t = 5.09; ES+: 838.36.

30 *(rac.)-(IR*, 5S*)-7-{4-[2-(4-Bromophenoxy)ethoxy]phenyl}-6-{cyclopropyl-[2-(2,3-difluorophenyl)ethyl]carbamoyl}-3,9-diazabicyclo[3.3.1]non-6-ene-9-carboxylic acid 2,2,2-trichloro-1,1-dimethylethyl ester hydrochloride salt (AL82)*

As for compound AL1 but from bicyclononene AK82 (0.26 mmol) in CH₂Cl₂ (5 mL) and HCl/dioxane (4M, 5 mL). LC-MS: R_t = 4.44; ES+: 840.42.

- 5 *(rac.)-(1R*, 5S*)-7-{4-[2-(4-Bromophenoxy)ethoxy]phenyl}-6-{cyclopropyl-[2-(4-fluorophenyl)ethyl]carbamoyl}-3,9-diazabicyclo[3.3.1]non-6-ene-9-carboxylic acid 2,2,2-trichloro-1,1-dimethylethyl ester hydrochloride salt (AL83)*

As for compound AL1 but from bicyclononene AK83 (0.24 mmol) in CH₂Cl₂ (5 mL) and HCl/dioxane (4M, 5 mL). LC-MS: R_t = 4.89; ES+: 822.39.

- 10 *(rac.)-(1R*, 5S*)-7-{4-[2-(4-Bromophenoxy)ethoxy]phenyl}-6-[cyclopropyl-(2-o-tolyethyl)carbamoyl]-3,9-diazabicyclo[3.3.1]non-6-ene-9-carboxylic acid 2,2,2-trichloro-1,1-dimethylethyl ester hydrochloride salt (AL84)*

- 15 As for compound AL1 but from bicyclononene AK84 (0.27 mmol) in CH₂Cl₂ (5 mL) and HCl/dioxane (4M, 5 mL).

- 20 *(rac.)-(1R*, 5S*)-7-{4-[2-(4-Bromophenoxy)ethoxy]phenyl}-6-[cyclopropyl-(3,5-dimethoxybenzyl)carbamoyl]-3,9-diazabicyclo[3.3.1]non-6-ene-9-carboxylic acid 2,2,2-trichloro-1,1-dimethylethyl ester hydrochloride salt (AL85)*

As for compound AL1 but from bicyclononene AK85 (0.25 mmol) in CH₂Cl₂ (5 mL) and HCl/dioxane (4M, 5 mL). LC-MS: R_t = 4.76; ES+: 850.40.

- 25 *(rac.)-(1R*, 5S*)-7-{4-[2-(4-Bromophenoxy)ethoxy]phenyl}-6-[cyclopropyl-(2-p-tolyethyl)carbamoyl]-3,9-diazabicyclo[3.3.1]non-6-ene-9-carboxylic acid 2,2,2-trichloro-1,1-dimethylethyl ester hydrochloride salt (AL86)*

- 30 As for compound AL1 but from bicyclononene AK86 (0.27 mmol) in CH₂Cl₂ (5 mL) and HCl/dioxane (4M, 5 mL). LC-MS: R_t = 4.99; ES+: 820.78.

(rac.)-(1R, 5S*)-7-{4-[2-(4-Bromophenoxy)ethoxy]phenyl}-6-{cyclopropyl-[2-(2-hydroxyethyl)benzyl]carbamoyl}-3,9-diazabicyclo[3.3.1]non-6-ene-9-carboxylic acid 2,2,2-trichloro-1,1-dimethylethyl ester hydrochloride salt (AL87)*

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As for compound **AL1** but from bicyclononene **AK87** (0.18 mmol) in CH₂Cl₂ (5 mL) and HCl/dioxane (4M, 5 mL). LC-MS: R_t = 4.58; ES+: 834.43.

(rac.)-(1R, 5S*)-7-Hydroxy-9-methyl-3,9-diazabicyclo[3.3.1]non-6-ene-3,6-dicarboxylic acid 6-benzyl ester 3-tert-butyl ester (AM)*

Ti(OEt)₄ (2.92 mL, 13.9 mmol) was added to a sol. of bicyclononane **A** (13.0 g, 39.8 mmol) in benzyl alcohol (90 mL). The mixture was heated to 125 °C and stirred at this temperature for 28 h. The mixture was allowed to cool to rt and aq. 10% HCl (180 mL) was added. The mixture was extracted with Et₂O (3x). The combined org. extracts were washed with aq. NaHCO₃ (2x), with brine (1x). The org. extracts were then dried over MgSO₄, filtered, and the solvents were removed first under reduced pressure, then under high vacuum. Purification of the residue by FC (EtOAc/heptane 1:1 → 3:1 → EtOAc) yielded the title compound (9.90 g, 64%). LC-MS: R_t = 1.39; ES+: 389.25.

(rac.)-(1R, 5S*)-9-Methyl-7-trifluoromethanesulfonyloxy-3,9-diazabicyclo[3.3.1]non-6-ene-3,6-dicarboxylic acid 6-benzyl ester 3-tert-butyl ester (AN)*

NaH (55% in oil, 2.20 g, 50.5 mmol) was added to a sol. of bicyclononane **AM** (15.69 g, 40.4 mmol) in THF (290 mL) at 0°C. After 15 min. Tf₂NPh (19.2 g, 53.7 mmol) was added and the mixture was stirred overnight while warming up to rt. Ice was added and the mixture was diluted with EtOAc, and washed with aq. 10% Na₂CO₃. The org. extracts were dried over MgSO₄, filtered, and the solvents were removed under reduced pressure. Purification of the residue by FC (EtOAc/heptane 1:1 → 3:1) yielded the title compound (17.1 g, 81%). R_f = 0.15 (EtOAc/heptane 1:1). LC-MS: R_t = 5.62; ES+: 521.37.

Compounds of type AO

(*rac.*)-(IR*, 5S*)-7-{4-[3-(*tert*-Butyldimethylsilyloxy)propyl]phenyl}-9-methyl-3,9-diazabicyclo[3.3.1]non-6-ene-3,6-dicarboxylic acid 6-benzyl ester 3-*tert*-butyl ester (AO1)

BuLi (1.5M in hexane, 3.81 mL, 5.71 mmol) was added to a sol. of [3-(4-bromophenyl)propoxy]-*tert*-butyldimethylsilane (Kiesewetter D. O., *Tetrahedron Asymmetry*, 1993, 4, 2183, 1.88 g, 5.71 mmol) in THF (33 mL) at -78 °C. After 30 min ZnCl₂ (1M in THF, 6.97 mL, 6.97 mmol, prepared as described for compound G1) was added and the mixture was allowed to warm up to rt. Bicyclononene AN (1.65 g, 3.17 mmol) and Pd(PPh₃)₄ (92 mg, 0.080 mmol) were added. The mixture was heated to 40 °C and stirred at this temperature for 30 min. The mixture was allowed to cool to rt and aq. 1M HCl (1 mL) was added. The mixture was diluted with EtOAc and washed with aq. 1M NaOH (1x). The org. extracts were dried over MgSO₄, filtered, and the solvents were removed under reduced pressure. Purification of the residue by FC (MeOH/CH₂Cl₂ 1:49 → 3:97 → 2:48 → 5:95) yielded the title compound (1.78 g, 90%). LC-MS: R_t = 5.55; ES+: 681.30.

(*rac.*)-(IR*, 5S*)-9-Methyl-7-{4-[2-(2,3,5-trimethylphenoxy)ethyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-3,6-dicarboxylic acid 6-benzyl ester 3-*tert*-butyl ester (AO2)

BuLi (1.6M in hexane, 19.40 mL, 31.0 mmol) was added to a sol. of compound C2 (9.90 g, 31.0 mmol) in THF (100 mL) at -78 °C. After 30 min ZnCl₂ (0.83M in THF, 43.8 mL, 37.2 mmol, prepared as described for compound G1) was added and the mixture was allowed to warm up to rt. Bicyclononene AN (9.90 g, 19.0 mmol) and Pd(PPh₃)₄ (550 mg, 0.475 mmol) were added. The mixture was heated to reflux for 1 h. The mixture was allowed to cool to rt and aq. 1M HCl (1 mL) was added. The mixture was diluted with EtOAc and washed with aq. 1M NaOH

(1x). The org. extracts were dried over MgSO₄, filtered, and the solvents were removed under reduced pressure. Purification of the residue by FC (MeOH/CH₂Cl₂ 1:49 → 3:97 → 2:48 → 5:95) yielded the title compound (6.20 g, 54%). LC-MS: R_t = 5.10; ES+: 611.59.

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Compounds of type AP

(*rac.*)-(1*R*^{*}, 5*S*^{*})-7-{4-[3-(*tert*-Butyldimethylsilyloxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-3,6,9-tricarboxylic acid 6-benzyl ester 3-*tert*-butyl ester 9-(2,2,2-trichloro-1,1-dimethylethyl) ester (AP1)

A sol. of bicyclononene AO1 (1.78 g, 2.87 mmol) and 2,2,2-trichloro-*tert*-butyl chloroformate (3.44 g, 14.4 mmol) in CH₂ClCH₂Cl (35 mL) was heated to reflux for 2 h. The mixture was allowed to cool to rt and the solvents were removed under reduced pressure. Purification of the residue by FC (EtOAc/heptane 1:8 → 1:1) yielded the title compound (1.88 g, 81%). LC-MS: R_t = 8.34.

(*rac.*)-(1*R*^{*}, 5*S*^{*})-7-{4-[2-(2,3,5-Trimethylphenoxy)ethyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-3,6,9-tricarboxylic acid 6-benzyl ester 3-*tert*-butyl ester 9-(2,2,2-trichloro-1,1-dimethylethyl) ester (AP2)

As for compound AP1 but from bicyclononene AO2 (22.4 g, 36.7 mmol) and 2,2,2-trichloro-*tert*-butyl chloroformate (44 g, 184 mmol) in CH₂ClCH₂Cl (400 mL). Purification of the residue by FC (EtOAc/heptane 1:8 → 1:1) yielded the title compound (19.2 g, 65%). LC-MS: R_t = 7.95.

(*rac.*)-(1*R*^{*}, 5*S*^{*})-3-Acetyl-7-[4-(3-hydroxypropyl)phenyl]-3,9-diazabicyclo[3.3.1]non-6-ene-6,9-dicarboxylic acid 6-benzyl ester 9-(2,2,2-trichloro-1,1-dimethylethyl) ester (AQ)

HCl/dioxane (4M, 20 mL) was added to a sol. of bicyclononene **AP1** (1.88 g, 2.32 mmol) in CH₂Cl₂ (20 mL) was cooled to 0 °C. The ice bath was removed and the mixture was stirred for 3 h at rt. The solvents were removed under reduced pressure and the residue was dried under high vacuum. This residue was then 5 dissolved in THF (30 mL) and the sol. was cooled to -78 °C. DMAP (cat. amount), DIPEA (1.60 mL, 9.28 mmol) and AcCl (0.165 mL, 2.32 mmol) were added. The mixture was stirred for 15 min at -78 °C and MeOH(10 mL) was added. The mixture was allowed to warm up to rt, was dissolved in EtOAc and washed with aq. 1M HCl (1x) and aq. sat. NaHCO₃ (1x). The org. Extracts were 10 dried over MgSO₄, filtered, and the solvents were removed under reduced pressure. Purification of the residue by FC (EtOAc/heptane 1...4 → 1:1 → EtOAc → MeOH/EtOAc 1:9) yielded the title compound (936 mg, 63%). LC-MS: R_t = 5.47; ES+: 637.06.

15 Compounds of type AR

(*rac.*)-(1*R*^{*, 5*S*^{*})-3-Acetyl-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6,9-dicarboxylic acid 6-benzyl ester 9-(2,2,2-trichloro-1,1-dimethylethyl) ester (**AR1**)}

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A mixture of bicyclononene **AQ** (468 mg, 0.73 mmol), 2,3,6-trifluorophenol (216 mg, 1.46 mmol), azodicarboxylic dipiperidine (277 mg, 1.10 mmol) and tributyl phosphine (0.541 mL, 2.19 mmol) in toluene (15 mL) was heated to reflux for 20 h. The mixture was allowed to cool to rt and the solvents were removed under 25 reduced pressure. Purification of the residue by FC (EtOAc/heptane 1:49 → 1:19 → 1:9) yielded the title compound (297 mg, 53%). LC-MS: R_t = 6.87; ES+: 767.04.

(*rac.*)-(1*R*^{*, 5*S*^{*})-3-Acetyl-7-{4-[3-(2-bromo-5-fluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6,9-dicarboxylic acid 6-benzyl ester 9-(2,2,2-trichloro-1,1-dimethylethyl) ester (**AR2**)}

As for the bicyclononene **AR1**, but from bicyclononene **AQ1** (468 mg, 0.73 mmol), 2-bromo-5-fluorophenol (0.163 mL, 1.46 mmol), azodicarboxylic dipiperidine (277 mg, 1.10 mmol) and tributyl phosphine (0.541 mL, 2.19 mmol) in toluene (15 mL). Purification of the residue by FC (EtOAc/heptane 1:49 → 5 1:19 → 1:9) yielded the title compound (205 mg, 35%). LC-MS: R_t = 7.06.

(*rac.*)-(1*R*^{*, 5*S*^{*})-3-Acetyl-6-[(2-allylbenzyl)cyclopropylcarbamoyl]-7-{4-[3-(2-bromo-5-fluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-9-carboxylic acid 2,2,2-trichloro-1,1-dimethylethyl ester (AS)}

10

A mixture of bicyclononene **AJ4** (225 mg, 0.300 mmol), (2-allylbenzyl)-cyclopropylamine (168 mg, 0.900 mmol), DIPEA (0.300 mL, 1.80 mmol), DMAP (10 mg, 0.082 mmol), HOEt (41 mg, 0.300 mmol) and EDC.HCl (86 mg, 0.450 mmol) in CH₂Cl₂ (3 mL) was stirred for 2 days. EDC.HCl (29 mg, 0.150 mmol) was added again after 24 h and 30 h. The mixture was diluted with CH₂Cl₂ and washed with aq. 1M HCl (1x). The org. extracts were dried over MgSO₄, filtered, and the solvents were removed under reduced pressure. Purification of the residue by FC (EtOAc/heptane 1:4 → 3:7 → 2:3 → 1:1 → 3:2 → 7:3 → 4:1) yielded the title compound (185 mg, 67%). R_f = 0.63 (EtOAc). LC-MS: R_t = 7.40.

15

1:1-Mixture of (*rac.*)-(1*R*^{*, 5*S*^{*})-3-acetyl-7-{4-[3-(2-bromo-5-fluorophenoxy)-propyl]phenyl}-6-{cyclopropyl-[2-((2*R*^{*)-2,3-dihydroxypropyl)benzyl]-carbamoyl}-3,9-diazabicyclo[3.3.1]non-6-ene-9-carboxylic acid 2,2,2-trichloro-1,1-dimethylethyl ester and (*rac.*)-(1*R*^{*, 5*S*^{*})-3-acetyl-7-{4-[3-(2-bromo-5-fluorophenoxy)propyl]phenyl}-6-{cyclopropyl-[2-((2*S*^{*)-2,3-dihydroxypropyl)benzyl]carbamoyl}-3,9-diazabicyclo[3.3.1]non-6-ene-9-carboxylic acid 2,2,2-trichloro-1,1-dimethylethyl ester (AT)}}}}

A mixture of bicyclononene **AS** (281 mg, 0.316 mmol), NMO·H₂O (44.8 mg, 30 0.332 mmol), and OsO₄ (2.5% in *tert*-BuOH, 0.0396 mL, 0.00316 mmol) in THF (4 mL), *tert*-BuOH (2 mL) and water (1 mL) was stirred overnight. NMO·H₂O (10 mg, 0.074 mmol) and OsO₄ (0.010 mL, 0.008 mmol) were added again and

the mixture was stirred again for 3 h. The solvents were removed under reduced pressure, and the residue was diluted with EtOAc, washed with aq. 1M HCl (1x), and with aq. sat. NaHCO₃ (1x). The org. extracts were dried over MgSO₄, filtered, and the solvents were removed under reduced pressure. Purification of 5 the residue by FC (EtOAc/heptane 1:4 → 2:3 → 3:2 → 4:1 → EtOAc → MeOH/EtOAc 1:9) yielded the title compounds (207 mg, 71%). R_f = 0.20 (EtOAc). LC-MS: R_t = 6.23; ES+: 922.59.

(*rac.*)-(1*R*<sup>*, 5*S*^{*})-3-Acetyl-7-{4-[3-(2-bromo-5-fluorophenoxy)propyl]phenyl}-
10 6-{cyclopropyl-[2-(2-oxoethyl)benzyl]carbamoyl}-3,9-diazabicyclo[3.3.1]non-
6-ene-9-carboxylic acid 2,2,2-trichloro-1,1-dimethylethyl ester (AU)</sup>

A mixture of bicyclononenes AT (167 mg, 0.181 mmol) and NaIO₄ (40 mg, 0.187 mmol) in THF (3 mL) and water (1 mL) was stirred at rt for 1 h. NaIO₄ (20 mg, 15 0.01 mmol) was added again and the mixture was stirred for 3 h. The mixture was diluted with EtOAc and washed with aq. sat. NaHCO₃ (1x). The org. extracts were dried over MgSO₄, filtered, and the solvents were removed under reduced pressure. The residue was dried under high vacuum and the title compound (156 mg, 97%) was used without further purification. LC-MS: R_t = 6.87; ES+: 891.78.

20 (rac.)-(1*R*<sup>*, 5*S*^{*})-3-Acetyl-7-{4-[3-(2-bromo-5-fluorophenoxy)propyl]phenyl}-
6-{cyclopropyl-[2-(2-hydroxyethyl)benzyl]carbamoyl}-3,9-
diazabicyclo[3.3.1]-non-6-ene-9-carboxylic acid 2,2,2-trichloro-1,1-
dimethylethyl ester (AV)</sup>

25 A mixture of bicyclononene AU (44.6 mg, 0.05 mmol) and NaBH₄ (about 2 mg, about 0.05 mmol) in MeOH (1 mL) was stirred at rt for 90 min. The mixture was diluted with EtOAc and washed with aq. 1M HCl (1x). The org. extracts were dried over MgSO₄, filtered and the solvents were removed under reduced 30 pressure. The residue was used without further purification.

Compounds of type AW

(*rac.*)-(1*R*^{*, 5*S*^{*})-7-[4-(3-Hydroxypropyl)phenyl]-3,9-diazabicyclo[3.3.1]non-6-ene-3,6,9-tricarboxylic acid 3-*tert*-butyl ester 6-ethyl ester 9-(2,2,2-trichloro-1,1-dimethylethyl) ester (AW1)}}

5

TBAF (28.8 g, 91.4 mmol) was added to a sol. of bicyclononene H3 (45.6 g, 60.9 mmol) in THF (900 mL) at 0 °C. After 20 min, the ice bath was removed. After stirring the mixture at rt for 5 h, it was diluted with EtOAc and washed with water (2x). The org. extracts were dried over MgSO₄, filtered, and the solvents were 10 removed under reduced pressure. Purification of the residue by FC (EtOAc/heptane 1:4 → 1:1) yielded the title compound (27.6 g, 72%). R_f = 0.22 (EtOAc/heptane 1:1). LC-MS: R_t = 6.11; ES+: 655.23.

(*rac.*)-(1*R*^{*, 5*S*^{*})-7-[4-(2-Hydroxyethoxy)phenyl]-3,9-diazabicyclo[3.3.1]non-15 6-ene-3,6,9-tricarboxylic acid 3-*tert*-butyl ester 6-ethyl ester 9-(2,2,2-trichloro-1,1-dimethylethyl) ester (AW2)}}

As for compound AW1 but from bicyclononene H7 (44.4 g, 59.2 mmol) TBAF (28.0 g, 88.9 mmol) and THF (600 mL). Purification by FC (EtOAc/heptane 1:3 20 → 1:1 → EtOAc) yielded the title compound (23.67 g, 63%). R_f = 0.20 (EtOAc/heptane 1:1). LC-MS: R_t = 6.02; ES+: 635.36.

Compounds of type AX

25 (*rac.*)-(1*R*^{*, 5*S*^{*})-7-{4-[3-(2,3,6-Trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-3,6,9-tricarboxylic acid 3-*tert*-butyl ester 6-ethyl ester 9-(2,2,2-trichloro-1,1-dimethylethyl) ester (AX1)}}

A mixture of bicyclononene AW1 (20.22 g, 32.0 mmol), 2,3,6-trifluorophenol 30 (9.50 g, 64.0 mmol), azodicarboxylic dipiperidine (16.15 g, 64.0 mmol) and tributyl phosphine (85%, 27.9 mL, 96.0 mmol) in toluene (800 ml) was heated to reflux for 2 h. The mixture was allowed to cool to rt and the solvent removed

under reduced pressure. Purification of the residue was purified by FC (EtOAc/heptane 1:19 → 1:9 → 1:4) yielded the title compound (21.7 g, 89%). $R_f = 0.60$ (EtOAc/heptane 1:1). LC-MS: $R_t = 1.25$; ES+: 765.22.

- 5 *(rac.)-(1R*, 5S*)-7-{4-[3-(2-Bromo-5-fluorophenoxy)propyl]phenyl}-3,9-diaza-bicyclo[3.3.1]non-6-ene-3,6,9-tricarboxylic acid 3-tert-butyl ester 6-ethyl ester 9-(2,2,2-trichloro-1,1-dimethylethyl) ester (AX2)*

As for AX1, but from bicyclononene AW1 (27.63 g, 43.6 mmol), 2-bromo-5-fluorophenol (9.70 mL, 87.2 mmol), azodicarboxylic dipiperidine (22.0 g, 87.2 mmol), tributyl phosphine (32.2 mL, 131 mmol), and toluene (550 mL). Purification of the residue by FC (EtOAc/heptane 1:19 → 1:9 → 1:4) yielded the title compound (31.67 g, 90%). $R_f = 0.60$ (EtOAc/heptane 1:1). LC-MS: $R_t = 7.63$.

- 15 *(rac.)-(1R*, 5S*)-7-{4-[2-(2-Chloro-4,5-dimethylphenoxy)ethoxy]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-3,6,9-tricarboxylic acid 3-tert-butyl ester 6-ethyl ester 9-(2,2,2-trichloro-1,1-dimethylethyl) ester (AX3)*

20 As for AX1, but from bicyclononene AW2 (11.83 g, 18.6 mmol), 2-chloro-4,5-dimethylphenol (5.83 mL, 37.2 mmol), azodicarboxylic dipiperidine (9.39 g, 37.2 mmol), tributyl phosphine (85%, 16.2 mL, 55.8 mmol), and toluene (300 mL). Purification of the residue by FC (EtOAc/heptane 1:19 → 1:9 → 1:3 → 1:1) yielded the title compound (13.35 g, 93%). $R_f = 0.50$ (EtOAc/heptane 1:1). LC-
25 MS: $R_t = 7.60$.

- (rac.)-(1R*, 5S*)-7-{4-[2-(4-Bromophenoxy)ethoxy]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6,9-dicarboxylic acid 6-ethyl ester 9-(2,2,2-trichloro-1,1-dimethylethyl) ester (AX4)*

30 As for AX1, but from bicyclononene AW2 (11.83 g, 18.6 mmol), 4-bromophenol (6.43 mL, 37.2 mmol), azodicarboxylic dipiperidine (9.39 g, 37.2 mmol), tributyl

phosphine (85%, 16.2 mL, 55.8 mmol), and toluene (300 ml). Purification of the residue by FC (EtOAc/heptane 1:19 → 1:9 → 1:3 → 1:1) yielded the title compound (13.6 g, 92%). $R_f = 0.50$ (EtOAc/heptane 1:1). LC-MS: $R_t = 7.49$.

5 Compounds of type AY

(*rac.*)-(1*R*^{*,} 5*S*^{*})-7-{4-[3-(2,3,6-Trifluorophenoxy)propyl]phenyl}-3,9-diaza-bicyclo[3.3.1]non-6-ene-3,6,9-tricarboxylic acid 3-*tert*-butyl ester 9-(2,2,2-trichloro-1,1-dimethylethyl) ester (AY1)

10

A mixture of bicyclononene AX1 (15.76 g, 20.7 mmol) in EtOH (600 mL) and aq. 1M NaOH (600 mL) was stirred for 7 h at 80 °C. The mixture was allowed to cool to rt and the solvents were partially removed under reduced pressure. The residue was diluted with EtOAc and aq. 1M HCl was added to pH 1 - 2. The 15 phases were shaken, separated and the aq. phase was extracted with EtOAc (2x). The combined org. extracts were dried over MgSO₄, filtered, and the solvents were removed under reduced pressure. Purification of the residue by FC (EtOAc/heptane 1:4 → 1:2 → 1:1) yielded the title compound (12.15 g, 80%). LC-MS: $R_t = 1.16$; ES+: 737.21.

20

(*rac.*)-(1*R*^{*,} 5*S*^{*})-7-{4-[3-(2-Bromo-5-fluorophenoxy)propyl]phenyl}-3,9-diaza-bicyclo[3.3.1]non-6-ene-3,6,9-tricarboxylic acid 3-*tert*-butyl ester 9-(2,2,2-trichloro-1,1-dimethylethyl) ester (AY2)

25

As for compound AY1, but from bicyclononene AX2 (29.67 g, 36.8 mmol), EtOH (700 mL) and aq. 1M NaOH (700 mL). Purification by FC (EtOAc/heptane 1:1 → EtOAc → MeOH/EtOAc 1:9) yielded the title compound 26.67 g (94%). LC-MS: $R_t = 6.89$; ES+: 749.92.

30

(*rac.*)-(1*R*^{*,} 5*S*^{*})-7-{4-[2-(2,3,5-Trimethylphenoxy)ethyl]phenyl}-3,9-diaza-bicyclo[3.3.1]non-6-ene-3,6,9-tricarboxylic acid 3-*tert*-butyl ester 9-(2,2,2-trichloro-1,1-dimethylethyl) ester (AY3)

Under N₂ Pd/C (10%, 1.92 g) was added to a sol. of bicyclononene AP2 (19.2 g, 24.0 mmol) in MeOH (390 mL) cooled to 0 °C. The mixture was purged with H₂ (4x) and stirred at 0 °C under H₂ for 7 h. The mixture was filtered through *Celite*,
5 diluted with EtOAc and washed with aq. 1M HCl (1x). The org. extracts were dried over MgSO₄, filtered, and the solvents were removed under reduced pressure. Purification by FC (EtOAc/heptane 1:3 → 1:1 → EtOAc → MeOH/EtOAc 1:9) yielded the title compound (4.26 g, 25%). LC-MS: R_t = 7.10.

10 (*rac.*)-(1*R*^{*,} 5*S*^{*})-7-{4-[2-(2-Chloro-4,5-dimethylphenoxy)ethoxy]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-3,6,9-tricarboxylic acid 3-*tert*-butyl ester 9-(2,2,2-trichloro-1,1-dimethylethyl) ester (AY4)

15 As for compound AY1, but from bicyclononene AX3 (13.35 g, 17.2 mmol), EtOH
→ EtOAc → MeOH/EtOAc 1:9) yielded the title compound 12.3 g (96%). R_f =
0.75 (EtOAc). LC-MS: R_t = 6.94.

20 (*rac.*)-(1*R*^{*,} 5*S*^{*})-7-{4-[2-(4-Bromophenoxy)ethoxy]phenyl}-3,9-diazabicyclo-[3.3.1]non-6-ene-6,9-dicarboxylic acid 9-(2,2,2-trichloro-1,1-dimethylethyl) ester (AY5)

25 As for compound AY1, but from bicyclononene AX4 (13.6 g, 17.2 mmol), EtOH
→ EtOAc → MeOH/EtOAc 1:9) yielded the title compound 12.2 g (93%). R_f =
0.75 (EtOAc). LC-MS: R_t = 6.75.

Compounds of type AZ

30 (*rac.*)-(1*R*^{*,} 5*S*^{*})-3-Acetyl-7-[4-(3-hydroxypropyl)phenyl]-3,9-diazabicyclo-[3.3.1]non-6-ene-6,9-dicarboxylic acid 9-(2,2,2-trichloro-1,1-dimethylethyl) ester (AZ1)

A mixture of bicyclononene S5 (3.23 g, 5.60 mmol) in EtOH (50 mL) and aq. 1M NaOH (50 mL) was stirred at 80 °C for 5 h. The mixture was allowed to cool to rt and diluted with EtOAc. The mixture was brought to pH 2 with aq. 1M HCl and 5 extracted with EtOAc (3x). The combined org. extracts were dried over MgSO₄, filtered, and the solvents were removed under reduced pressure. Purification by FC (MeOH/CH₂Cl₂ 1:19 → 1:9 → 1:4) yielded the title compound (1.40 g, 46%). LC-MS: R_t = 0.89; ES+: 547.28.

10 (*rac.*)-(1*R*^{*,} 5*S*^{*)}-7-[4-(2-Hydroxyethyl)phenyl]-3,9-diazabicyclo[3.3.1]non-6-ene-3,6,9-tricarboxylic acid 3-*tert*-butyl ester 9-(2,2,2-trichloro-1,1-dimethyl-ethyl) ester (AZ2)

15 As for compound AZ1 but from bicyclononene H8 (4.96 g), EtOH (150 mL) and aq. 1M NaOH (150 mL). The crude material was used further without purification.

Compounds of type BA

20 (*rac.*)-(1*R*^{*,} 5*S*^{*)}-3-Acetyl-7-{4-[3-(*tert*-butyldimethylsilyloxy)propyl]-phenyl}-6-[(2-chlorobenzyl)cyclopropylcarbamoyl]-3,9-diazabicyclo[3.3.1]-non-6-ene-9-carboxylic acid 2,2,2-trichloro-1,1-dimethylethyl ester (BA1)

25 A mixture of bicyclononene T4 (1.85 g, 2.79 mmol), (2-chlorobenzyl)-cyclopropylamine (1.52 g, 8.37 mmol), DMAP (85 mg, 0.70 mmol), DIPEA (1.91 mL, 11.2 mmol), HOBr (377 mg, 2.79 mmol) and EDC·HCl (803 mg, 4.19 mmol) in CH₂Cl₂ (50 mL) was stirred at rt for 18 h. The mixture was diluted with more CH₂Cl₂ and washed with aq. 1M HCl (1x) and aq. sat. NaHCO₃ (1x). The org. extracts were dried over MgSO₄, filtered, and the solvents were removed under 30 reduced pressure. Purification of the residue by FC (EtOAc/heptane 1:4 → 1:3 → 1:2 → 1:1) yielded the title compound (1.16 g, 50%). LC-MS: R_t = 1.37.

(*rac.*)-(1*R*^{*,}, 5*S*^{*)}-7-{4-[2-(*tert*-Butyldiphenylsilanyloxy)ethyl]phenyl}-6-[(2-chlorobenzyl)cyclopropylcarbamoyl]-3,9-diazabicyclo[3.3.1]non-6-ene-3,9-dicarboxylic acid 3-*tert*-butyl ester 9-(2,2,2-trichloro-1,1-dimethylethyl) ester (BA2)

5

As for compound BA1, but from bicyclononene T5 (crude, about 5.79 mmol), (2-chlorobenzyl)cyclopropylamine (3.10 g, 17.1 mmol), DIPEA (3.9 mL, 22.8 mmol), DMAP (140 mg, 1.14 mmol), HOBr (770 mg, 5.70 mmol), and EDC-HCl (1.64 g, 8.55 mmol), in CH₂Cl₂ (50 mL). Purification of the residue by FC (EtOAc/heptane 1:8 → 1:4) yielded the title compound 3.35 g (58%). R_f = 0.55 (EtOAc/heptane 2:3). LC-MS: R_t = 1.40.

(*rac.*)-(1*R*^{*,}, 5*S*^{*)}-3-Acetyl-7-{4-[2-(*tert*-butyldiphenylsilanyloxy)ethyl]phenyl}-6-[(2-chlorobenzyl)cyclopropylcarbamoyl]-3,9-diazabicyclo[3.3.1]non-6-ene-9-carboxylic acid 2,2,2-trichloro-1,1-dimethylethyl ester (BA3)

As for compound S1, from BA2 (1.45 g, 1.45 mmol), CH₂Cl₂ (10 mL), 4M HCl/dioxane (10 mL), THF (20 mL), without DMAP, DIPEA (4.62 mL, 27.0 mmol), acetyl chloride (0.903 mL, 9.55 mmol), and MeOH (5 mL). Purification of the residue by FC (ErOAc/heptane 1:1 → EtOAc → MeOH/EtOAc 1:9) led to the title compound (1.34 g, 75%). R_f = 0.30 (EtOAc/heptane 1:1). LC-MS: R_t = 1.39.

Compounds of type BB

25

(*rac.*)-(1*R*^{*,}, 5*S*^{*)}-3-Acetyl-6-[(2-chlorobenzyl)cyclopropylcarbamoyl]-7-{4-(3-hydroxypropyl)phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-9-carboxylic acid 2,2,2-trichloro-1,1-dimethylethyl ester (BB1)

30 A mixture of bicyclononene BA1 (1.16 g, 1.40 mmol) and TBAF (884 mg, 2.80 mmol) in THF (10 mL) was stirred at rt for 90 min. The mixture was diluted with EtOAc and washed with water (2x) and brine (1x). The org. extracts were dried

over MgSO₄, filtered, and the solvents were removed under reduced pressure. Purification by FC (MeOH/CH₂Cl₂ 1:49 → 1:9) yielded the title compound (990 mg, 98%). R_f = 0.47 (MeOH/CH₂Cl₂ 1:9). LC-MS: R_t = 1.11.

- 5 **(rac.)-(IR*, 5S*)-3-Acetyl-6-[(2-chlorobenzyl)cyclopropylcarbamoyl]-7-[4-(2-hydroxyethyl)phenyl]-3,9-diazabicyclo[3.3.1]non-6-ene-9-carboxylic acid 2,2,2-trichloro-1,1-dimethylethyl ester (BB2)**

As for compound BB1, but from BA3 (1.99 g, 2.21 mmol), TBAF (1M in THF, 10 4.5 mL, 4.5 mmol) in THF (15 mL). Purification by FC (EtOAc/heptane 1:5 → 1:1 → EtOAc) yielded the title compound (1.00 g, 68%). R_f = 0.38 (EtOAc/heptane 1:1). LC-MS: R_t = 1.09; ES+: 698.02.

- 15 **(rac.)-(IR*, 5S*)-6-[(2-Chlorobenzyl)cyclopropylcarbamoyl]-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-3-carboxylic acid *tert*-butyl ester (BC)**

Zn (1.63 g, 24.9 mmol) was added to a sol. of bicyclononene AK2 (2.25 g, 2.50 mmol) in THF (30 mL) and AcOH (10 mL) under efficient stirring. The mixture 20 was stirred efficiently for 2.5 h, then filtered and washed with THF. The filtrate was diluted with EtOAc and washed with aq. 1M NaHO (2x). The org. extracts were dried over MgSO₄, and filtered. Evaporating the solvents under reduced pressure yielded the title compound that was used without further purification.

- 25 **(IR, 5S)-9-Methyl-7-trifluoromethanesulfonyloxy-3,9-diazabicyclo[3.3.1]non-6-ene-3,6-dicarboxylic acid 6-benzyl ester 3-*tert*-butyl ester (BD)**

A sol. of bicyclononene AG (1.13 g; 2.91 mmol) in THF (8 ml) was added to a suspension of NaH (ca 60%, 175 mg; 4.36 mmol) in THF (2 ml) at 0°C. After 30 30 min Tf₂NPh (1.56 g; 4.36 mmol) was added and the mixture was stirred at rt for 12 h. Ice (5 g) was added and THF was evaporated. The aq. residue was extracted with EtOAc (3x). The combine org. extracts were dried over MgSO₄, filtered, and

the solvents were removed under reduced pressure. Purification of the residue by FC (EtOAc/cyclohexane 1:1 → EtOAc) yielded the title compound (1.28 g, 84%). R_f = 0.53 (EtOAc).

5 (*1R, 5S*)-7-{4-[3-(*tert*-Butyldimethylsilyloxy)propyl]phenyl}-9-methyl-3,9-diazabicyclo[3.3.1]non-6-ene-3,6-dicarboxylic acid 6-benzyl ester 3-*tert*-butyl ester (BE)

BuLi (1.6 M in hexane, 3.82 mL, 5.98 mmol) was added to a sol. of [3-(4-bromophenyl)propoxy]-*tert*-butyldimethylsilane (Kiesewetter D. O., *Tetrahedron Asymmetry*, 1993, 4, 2183; 1.97 g; 5.98 mmol) in THF (4ml) at -78°C. After 30 min, ZnCl₂ (1M in THF, 7.2 ml; 7.2 mmol) was added and the mixture was allowed to warm up to rt. A sol. of bicyclononene BD (1.24 g; 2.39 mmol) in THF (7 ml) and Pd(PPh₃)₄ (69 mg; 0.060 mmol) were added after each other and the mixture was heated to 40°C for 35 min. The reaction mixture was allowed to cool to rt, sat. solution of NH₄Cl was added, and the mixture was extracted with EtOAc (3x). The combined org. extracts were dried over MgSO₄, filtered, and the solvents were removed under reduced pressure. Purification of the residue by FC (EtOAc/cyclohexane 1:1 → EtOAc) yielded the title compound (1.22 g, 82%). R_f = 0.27 (EtOAc). LC-MS: R_t = 5.68; ES+ = 621.30.

(*1R, 5S*)-7-{4-[3-(*tert*-Butyldimethylsilyloxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-3,6,9-tricarboxylic acid 6-benzyl ester 3-*tert*-butyl ester 9-(2,2,2-trichloro-1,1-dimethylethyl) ester (BF)

A mixture of bicyclononene BE (1.66 g, 2.67 mmol) and β,β,β-trichloro-*tert*-butyl chloroformate (13.4 g, 240 mmol) in 1,2-dichloroethane (10 ml) was heated to reflux for 4 h. The mixture was allowed to cool to rt and the solvents were removed under reduced pressure. Purification of the residue by FC (EtOAc/cyclohexane 1:4) yielded the title compound (1.75 g, 83%). R_f = 0.43 (EtOAc/cyclohexane 1:4). LC-MS: R_t = 8.30.

(*1R*, *5S*)-7-[4-(3-Hydroxypropyl)phenyl]-3,9-diazabicyclo[3.3.1]non-6-ene-3,6,9-tricarboxylic acid 3-*tert*-butyl ester 9-(2,2,2-trichloro-1,1-dimethylethyl) ester (BH)

5 A sol. of bicyclononene **BF** (1.59 g, 1.96 mmol) in CH₂Cl₂ (3 ml) was cooled to 0°C and HCl/dioxane (4M, 10 ml) was added. The mixture was stirred for 2 h at 0°C and subsequently for 3 h at rt. After the solvents were removed under reduced pressure the crude was dried under high vacuum. The cresidue was dissolved in THF (5 ml). DMAP (12 mg, 0.098 mmol) and DIPEA (1.34 ml; 10 7.849 mmol) were added and the mixture was cooled to -78 °C. AcCl (0.153 ml; 2.16 mmol) was added and reaction mixture was stirred at -78 °C for 30 min. After addition of MeOH (1 ml) and warming-up to rt, aq. HCl (1M, 10 ml) was added and reaction mixture was extracted with EtOAc (3x). The combined org. extracts were washed with aq. sat. NaHCO₃ (1x), and the org. extracts were dried 15 over MgSO₄, filtered, and the solvents were removed under reduced pressure. Purification of the residue by FC EtOAc/cyclohexane 1:3 → 1:1) yielded the title compound (280 mg, 22%). R_f = 0.38 (EtOAc). LC-MS: R_t = 5.43; ES+ = 637.17.

(*1R*, *5S*)-7-[4-(3-Hydroxypropyl)phenyl]-3,9-diazabicyclo[3.3.1]non-6-ene-20 3,6,9-tricarboxylic acid 3-*tert*-butyl ester 9-(2,2,2-trichloro-1,1-dimethylethyl) ester (BI)

A mixture of bicyclononene **BH** (140 mg; 0.219 mmol) and Pd/C (10%, 25 mg) in MeOH (4 ml) was stitted at rt under H₂ for 2 h. The mixture was filtered through 25 Celite, washed with MeOH, and the solvents were evaporated under reduced pressure. The crude product (110 mg) was directly used in the next reaction without purification. R_f = 0.15 (EtOAc). LC-MS: R_t = 4.41; ES- : 545.02.

(*1R*, *5S*)-7-[4-(3-Hydroxypropyl)phenyl]-6-(methylphenethylcarbamoyl)-3,9-30 diazabicyclo[3.3.1]non-6-ene-3,9-dicarboxylic acid 3-*tert*-butyl ester 9-(2,2,2-trichloro-1,1-dimethylethyl) ester (BJ)

A mixture of bicyclononene **BI** (95 mg; 0.17 mmol), phenethylmethylamine (0.48 ml; 0.34 mmol); HOEt (6.0 mg, 0.042 mmol), EDC·HCl (49 mg; 0.255 mmol) and DMAP (5.0 mg; 0.042 mmol) in CHCl₃ (6 ml) was stirred at rt for 14 h. Aq HCl (1M) was added and the mixture was extracted with CH₂Cl₂ (3x). The org. phase was washed with aq. sat. NaHCO₃ (1x), the combined org. extracts were dried over MgSO₄, filtered, and the solvents were removed under reduced pressure. Purification of the residue by FC (EtOAc/cyclohexane 1:1 → EtOAc) yielded the title compound (64 mg, 44%). R_f = 0.25 (EtOAc). LC-MS: R_t = 5.37; ES+:
5 664.29.

10

(IR, *5S)-7-{4-[3-(2-Bromo-5-fluorophenoxy)propyl}phenyl]-6-(methyl-phenethylcarbamoyl)-3,9-diazabicyclo[3.3.1]non-6-ene-3,9-dicarboxylic acid 3-*tert*-butyl ester 9-(2,2,2-trichloro-1,1-dimethyl-ethyl) ester (BK)*

15 A mixture of bicyclononene **BJ** (60 mg; 0.090 mmol), 2-bromo-5-fluorophenol
 (34 mg, 0.18 mmol), azodicarboxylic dipiperide (34 mg; 0.135 mmol) and
 tributylphosphine (67 mg; 0.270 mmol) in toluene (2 ml) was heated to reflux for
 20 h. The solvent was removed under reduced pressure. Purification of the
 residue was by FC EtOAc/cyclohexane 2:1 → 4:1 yielded the title compound (58
 mg, 76%). $R_f = 0.60$ (EtOAc). LC-MS: $R_t = 7.01$; ES+ = 836.07.

(*rac*.)-(1*R*^{*}, 5*S*^{*})-3-Acetyl-7-{4-[3-(2-bromo-5-fluorophenoxy)-propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6,9-dicarboxylic acid 6-ethyl ester 9-(2,2,2-trichloro-1,1-dimethylethyl) ester (**BL**)

25

HCl/dioxane (4M, 20 mL) was added to a sol. of bicyclononene **AX2** (2.00 g, 2.47 mmol) in CH₂Cl₂ (20 mL) cooled to 0 °C. The ice bath removed and the mixture was stirred at rt for 2 h. The solvents were removed under reduced pressure and the foamy residue dried under high vacuum. A mixture of this residue, DMAP (15 mg, 0.123 mmol) and DIPEA (1.69 mL, 9.88 mmol) in THF (40 mL) was cooled to -78 °C, and AcCl (0.186 mL, 2.47 mmol) was added. The mixture was stirred for 20 min at -78 °C and MeOH (5 mL) was added. The mixture was

allowed to warm up to rt, was diluted with EtOAc and washed with aq. 1M HCl (1x). The org. extracts were dried over MgSO₄, filtered, and the solvents were removed under reduced pressure. Purification of the residue by FC (EtOAc/heptane 1:4 → 1:1 → EtOAc) yielded the title compound (1.55 g, 83%).

5 R_f = 0.50 (EtOAc).

Compounds of type BM

10 3-Acetyl-7-{4-[2-(2-bromo-5-fluorophenoxy)ethoxy]phenyl}-3,9-diaza-bicyclo[3.3.1]non-6-ene-6,9-dicarboxylic acid 6-ethyl ester 9-(2,2,2-trichloro-1,1-dimethylethyl) ester (**BM2**)

15 Tributylphosphine (3.84 mL, 15.6 mmol) was added to a sol. of bicyclononene **S4** (3.00 g, 5.19 mmol), 2-bromo-5-fluorophenol (1.15 mL, 10.4 mmol) and azodicarboxylic dipiperidide (1.97 g, 7.79 mmol) in toluene (30 mL). The mixture was heated to reflux for 2 h and allowed to cool to rt. The solvents were removed under reduced pressure. Purification by FC (EtOAc/heptane 1:1 → 2:1 → 3:1) yielded the title compound (2.70 g, 69%).

20 3-Acetyl-7-{4-[2-(2-chloro-4,5-dimethylphenoxy)ethoxy]phenyl}-3,9-diaza-bicyclo[3.3.1]non-6-ene-6,9-dicarboxylic acid 6-ethyl ester 9-(2,2,2-trichloro-1,1-dimethylethyl) ester (**BM3**)

25 As described for compound **BM2**, but from bicyclononene **S4** (3.00 g, 5.19 mmol), 2-chloro-4,5-dimethylphenol (1.64 g, 10.5 mmol), azodicarboxylic dipiperidide (1.98 g, 7.86 mmol), tributylphosphine (3.90 mL, 15.7 mmol) and toluene (50 mL). Purification by FC yielded the title compound (2.82 g, 75%).

30 3-Acetyl-7-{4-[2-(2,6-dichloro-4-methylphenoxy)ethoxy]phenyl}-3,9-diaza-bicyclo[3.3.1]non-6-ene-6,9-dicarboxylic acid 6-ethyl ester 9-(2,2,2-trichloro-1,1-dimethylethyl) ester (**BM4**)

As described for compound **BM2**, but from bicyclononene **S4** (3.00 g, 5.19 mmol), 2,6-dichloro-4-methylphenol (1.84 g, 10.38 mmol), azodicarboxylic dipiperide (1.97 g, 7.79 mmol) , tributylphosphine (3.84 mL, 15.6 mmol) and toluene (50 mL). Purification by FC yielded the title compound (2.76 g, 72%).

5

3-Acetyl-7-{4-[2-(2,3-dichlorophenoxy)ethoxy]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6,9-dicarboxylic acid 6-ethyl ester 9-(2,2,2-trichloro-1,1-dimethylethyl) ester (BM5)

10 As described for compound **BM2**, but from bicyclononene **S4** (3.20 g, 5.54 mmol), 2,3-dichlorophenol (1.80 g, 11.1 mmol), azodicarboxylic dipiperide (2.10 g, 8.31 mmol) , tributylphosphine (4.11 mL, 16.6 mmol) and toluene (50 mL). Purification by FC yielded the title compound (2.22 g, 55%).

15 **3-Acetyl-7-{4-[2-(4-chloro-2-methylphenoxy)ethoxy]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6,9-dicarboxylic acid 6-ethyl ester 9-(2,2,2-trichloro-1,1-dimethylethyl) ester (BM7)**

20 As described for compound **BM2**, but from bicyclononene **S4** (3.00g, 5.19 mmol), 4-chloro-2-methylphenol (1.48 g, 10.4 mmol), azodicarboxylic dipiperide (1.97 g, 7.79 mmol) , tributylphosphine (3.84 mL, 15.6 mmol) and toluene (50 mL). Purification by FC yielded the title compound (1.36 g, 37%).

25 **3-Acetyl-7-{4-[2-(2,4,5-trichlorophenoxy)ethoxy]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6,9-dicarboxylic acid 6-ethyl ester 9-(2,2,2-trichloro-1,1-dimethylethyl) ester (BM9)**

30 As described for compound **BM2**, but from bicyclononene **S4** (3.00 g, 5.19 mmol) 2,4,5-trichlorophenol (2.05 g, 10.4 mmol), azodicarboxylic dipiperide (1.97 g, 7.79 mmol) , tributylphosphine (3.84 mL, 15.6 mmol) and toluene (50 mL). Purification by FC yielded the title compound (2.76 g, 72%).

3-Acetyl-7-{4-[2-(2-chloro-5-fluorophenoxy)ethoxy]phenyl}-3,9-diaza-bicyclo[3.3.1]non-6-ene-6,9-dicarboxylic acid 6-ethyl ester 9-(2,2,2-trichloro-1,1-dimethylethyl) ester (BM10)

5 As described for compound **BM2**, but from bicyclononene **S4** (3.18 g, 5.50 mmol) 2-chloro-5-fluorophenol (1.61 g, 11.0 mmol), azodicarboxylic dipiperide (2.08 g, 8.25 mmol), tributylphosphine (4.10 mL, 16.6 mmol) and toluene (50 mL). Purification by FC yielded the title compound (2.67 g, 69%).

10 **Compounds of type BN**

3-Acetyl-7-{4-[2-(2-bromo-5-fluorophenoxy)ethoxy]phenyl}-3,9-diaza-bicyclo[3.3.1]non-6-ene-6,9-dicarboxylic acid 9-(2,2,2-trichloro-1,1-dimethyl-ethyl) ester (BN2)

15 A mixture of bicyclononene **BM2** (2.69, 3.58 mmol) in aq. 1M NaOH (30 mL) and EtOH (70 mL) was stirred for 1 h at 85 °C. The mixture was allowed to cool to rt and the solvents were partially removed under reduced pressure. The residue was acidified to pH 2 with aq. 1M HCl and this mixture was extracted with EtOAc (3x). The combined org. extracts were dried over MgSO₄, filtered, and the solvents were removed under reduced pressure. The crude title compound (2.96 g, quantitative yield) was used further without purification.

20 **3-Acetyl-7-{4-[2-(2-chloro-4,5-dimethylphenoxy)ethoxy]phenyl}-3,9-diaza-bicyclo[3.3.1]non-6-ene-6,9-dicarboxylic acid 9-(2,2,2-trichloro-1,1-dimethyl-ethyl) ester (BN3)**

25 As described for compound **BN2**, but from bicyclononene **BM3** (2.82 g, 3.94 mmol), aq. 1M NaOH (30 mL) and EtOH (70 mL). The crude title compound (30 g, 96%) was used further without purification.

3-Acetyl-7-{4-[2-(2,6-dichloro-4-methylphenoxy)ethoxy]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6,9-dicarboxylic acid 9-(2,2,2-trichloro-1,1-dimethyl-ethyl) ester (BN4)

5 As described for compound BN2, but from bicyclononene BM4 (2.75 g, 3.73 mmol), aq. 1M NaOH (30 mL) and EtOH (70 mL). The crude title compound (2.63 g, quantitative yield) was used further without purification.

3-Acetyl-7-{4-[2-(2,3-dichlorophenoxy)ethoxy]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6,9-dicarboxylic acid 9-(2,2,2-trichloro-1,1-dimethylethyl) ester (BN5)

10 As described for compound BN2, but from bicyclononene BM5 (2.22 g, 3.07 mmol), aq. 1M NaOH (30 mL) and EtOH (70 mL). The crude title compound (1.59 g, 75%) was used further without purification.

3-Acetyl-7-{4-[2-(4-chloro-2-methylphenoxy)ethoxy]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6,9-dicarboxylic acid 9-(2,2,2-trichloro-1,1-dimethyl-ethyl) ester (BN7)

15 20 As described for compound BN2, but from bicyclononene BM7 (1.35 g, 1.92 mmol), aq. 1M NaOH (30 mL) and EtOH (70 mL). The crude title compound (1.25 g, 97%) was used further without purification.

3-Acetyl-7-{4-[2-(2,4,5-trichlorophenoxy)ethoxy]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6,9-dicarboxylic acid 9-(2,2,2-trichloro-1,1-dimethylethyl) ester (BN9)

25 As described for compound BN2, but from bicyclononene BM9 (2.31 g, 3.05 mmol), aq. 1M NaOH (30 mL) and EtOH (70 mL). The crude title compound (2.19 g, 99%) was used further without purification.

3-Acetyl-7-{4-[2-(2-chloro-5-fluorophenoxy)ethoxy]phenyl}-3,9-diazabicyclo-[3.3.1]non-6-ene-6,9-dicarboxylic acid 9-(2,2,2-trichloro-1,1-dimethylethyl)ester (BN10)

- 5 As described for compound BN2, but from bicyclononene BM10 (2.82 g, 3.94 mmol), aq. 1M NaOH (30 mL) and EtOH (70 mL). The crude title compound (1.90 g, 74%) was used further without purification.

Preparation of the final compounds

10

Typical procedure (A) for the acylation:

- To a solution of bicyclononene in anhydrous EtOAc was added vacuum dried-Amberlyst 21 (1.5 g/mmol of bicyclononene) or another suitable scavenger, 15 followed by the addition of the desired acid chloride (1.5 eq.). After shaking the suspension for 3 h, an aliquot water was added and shaking was continued for 1 h. The resin was then removed by filtration, washed with EtOAc and the filtrate was evaporated to yield the intermediate amide.

- 20 The synthesis of the sulfonamide, carbamate or urea derivatives was performed in analogy to the above-described procedure, by using the corresponding sulfonyl chloride, chloroformate or carbamoyl chloride respectively.

Typical procedure (B) for amide formation from acid chlorides:

25

- To a sol. of the acid chloride (1 eq.) in CH₂Cl₂ (2.5 mL/mmol) at 0 °C. the amine (3 eq.) was added. The mixture was stirred for 3 h while warming up slowly to rt. If necessary, more CH₂Cl₂ was added, then the reaction mixture was washed with aq. sat. NaHCO₃ (1x) and aq. 1M HCl (1x). The extracts were dried over MgSO₄, 30 and the solvents were removed under reduced pressure. The obtained product was used without further purification.

Typical procedure (C) for an amide coupling with CDI

To a sol. of the carboxylic acid (1 eq.) in CH₂Cl₂ (4 mL/mmol) CDI (1 eq.) was added. The sol. or suspension was stirred for 2 h at rt, then cooled to 0 °C. The 5 amine (6 eq.) was added and the sol. or suspension was stirred for 2 h while warming up slowly to rt. The sol. or suspension was washed with water (1x). The org. extracts were evaporated under reduced pressure and the obtained residue was used further without purification.

10 *Typical procedure (D) for the reduction of an amide to an amine with LAH*

To a sol. of the amide (1 eq.) was dissolved in THF (3 mL/mmol) LAH (1M in THF, 3 eq.) was added carefully. The mixture was stirred at rt for 30 min, then heated to 60 °C for 3 h before it was allowed to cool down to rt, then to 0 °C. For 15 x g of LiAlH₄ initially added, was added x g of water, then x g of aq. 15% NaOH, and finally 3x g of water again. The resulting mixture was stirred overnight, filtered, and the precipitate washed with EtOAc. The filtrate was evaporated under reduced pressure and the residue diluted in a small amount of MeOH. The sol. was passed through a pad of SCX silica gel (sulfonic acid). Elution started 20 with MeOH, followed by NH₃/MeOH. The amines eluted with the second second eluent. The solvents were removed under reduced pressure. The isolated amines were either used without further purification or purified by HPLC, depending on the purity.

25 *Typical procedure (E) for the cleavage of the 2,2,2-trichloro-1,1-dimethylethylcarbamate protecting group:*

The crude product from another typical procedure was dissolved in THF/AcOH (1:0.1) and treated with zinc (20 eq.). The suspension was stirred for 5 h and 30 filtered through celite, which was washed with EtOAc. The filtrate was evaporated under reduced pressure and the residue was purified by HPLC.

Typical procedure (F) for the formation of aryl ether (Mitsunobu reaction)

The bicyclononene (0.05 mmol) was dissolved or suspended in toluene (1.00 mL). The phenol derivative (0.075 mmol) in toluene (0.50 mL) was added. TMAD 5 (0.075 mmol) in toluene (0.50 mL) was added, followed by tributylphosphine (0.15 mmol). The reaction mixture was stirred for 2 h at rt and then 2 h at 60 °C. Sometimes, it was necessary to add a second portion of tributylphosphine and to stir overnight. Sometimes, THF was necessary as cosolvent to dissolve the reactants. The reaction mixture was allowed to cool to rt, then water was added. 10 The mixture was extracted with EtOAc, and the org. extracts were evaporated under reduced pressure.

Typical procedure (G) for an amide coupling

15 To a sol. of bicyclononene (0.05 mmol) in CHCl₃ (2 mL) the desired carboxylic acid (0.10 mmol) was added. DIPEA (0.10 mmol), DMAP (0.01 mmol), HOBr (0.01 mmol), and EDC·HCl (0.05 mmol) were added and the reaction mixture was stirred overnight. Sometimes, it was necessary to add another portion of acid, DMAP, HOBr and EDC·HCl and to continue stirring for 24 h. CH₂Cl₂ was added 20 and the mixture was washed with water. The org. extracts were separated, dried over Na₂SO₄ and filtered. The solvent was removed under reduced pressure.

Typical procedure (H) for an amide coupling

25 To a sol. of the bicyclononene (0.05 mmol) CHCl₃ or CH₂Cl₂ (2 mL) the desired amine (commercially available or prepared following known, standard procedures) (0.10 mmol) was added. DIPEA (0.10 mmol), DMAP (0.01 mmol), HOBr (0.01 mmol) and EDC·HCl (0.05 mmol) were added. The reaction mixture was stirred overnight. Sometimes, it was necessary to add another portion of 30 amine, DMAP, HOBr and EDC·HCl and to continue stirring the sol. for 24 h. CH₂Cl₂ was added and the mixture was washed with water. The org. extracts

were separated, dried over Na_2SO_4 and filtered. The solvents were removed under reduced pressure.

Typical procedure (J) for reductive amination

5

To a solution of aldehyde (1 eq.) in MeOH (0.5 mL/mmol) was added an amine (1.2 eq.). The solution was stirred for 2h. Sodium borohydride (1.2 eq.) was added portionwise at 0°C and then stirring was continued, at rt, for 4h. A solution of NaOH 1N was added and the MeOH was evaporated. The mixture was extracted 10 with EtOAc twice and the organic layer was washed with brine, dried over Na_2SO_4 and filtered. The solvent was removed under reduced pressure. The isolated amines were either used without further purification or purified by flash chromatography (EtOAc/heptane: 2/8), depending on the purity.

15 *Typical procedure (K) for an anhydride coupling*

To a sol. of the bicyclononene (0.05 mmol) in CH_2Cl_2 (0.4 mL) was added DIPEA (0.1 mmol) followed by the anhydride (0.06 mmol) in CH_2Cl_2 (0.4 mL) at 0°C. After stirring for 3h at rt, the solvent was evaporated under reduced pressure.

20

Typical procedure (L) for protecting group (BOC and TBDMS) cleavage

To a sol. of the bicyclononene (0.05 mmol) in CH_2Cl_2 (0.5 mL), cooled to 0°C, was added 4M HCl/dioxane(0.5 mL). The ice bath was removed and the solution 25 was stirred for 1h30 to 3h, depending on the compound. The solvents were evaporated under reduced pressure without heating.

Typical procedure (M) for the saponification of esters

30 A mixture of the ester (1 eq.) and LiOH (2 eq.) in THF was stirred at rt for 2 h. The solvents were removed under reduced pressure and the residue was extracted on isolute sorbent (0.25 g pre-washed with 0.300 mL aq. 1M HCl, elution with 2

mL CH₂Cl₂). The solvent was removed under reduced pressure and the residue was used without further purification.

Preparation of amines

5

(2-Chlorobenzyl)cyclopropylamine

Synthesized according to typical procedures B and D from 2-chlorobenzoyl chloride and cyclopropylamine.

10

Benzylcyclopropylamine

See Loepky, R. N.; *et al.*, *J. Org. Chem.*, **2000**, *65*, 96.

15 **(2-Chlorobenzyl)ethylamine**

See Ishihara, Y; *et al.*; *Chem. Pharm. Bull.*, **1991**, *39*, 3225.

Cyclopropyl-(3-trifluoromethylbenzyl)amine

20

See Brabander, H. J.; *et al.*; *J. Org. Chem.*, **1967**, *32*, 4053.

Cyclopropylphenethylamine

25

See Smith, P. W.; *et al.*; *J. Med. Chem.*, **1998**, *41*, 787.

Methyl(3-phenoxypropyl)amine

30

Synthesized according to typical procedures C and D from 3-phenoxypropionic acid and methylamine.

(2-p-Tolyloxyethyl)methylamine

Synthesized according to typical procedures C and D from 2-*p*-tolyloxyacetic acid and methylamine.

5 [2-(3-Chlorophenyl)ethyl]amine

Synthesized according to typical procedures C and D from 3-chlorophenylacetic acid and methylamine.

10 [2-(2-Methoxyphenyl)ethyl]amine

Synthesized according to typical procedures C and D from 2-methoxyphenylacetic acid and methylamine.

15 (2-Allylbenzyl)cyclopropylamine

BuLi (1.55 M in hexane, 14.7 mL, 22.7 mmol) was added to a sol. of 1-bromo-2-(dimethoxymethyl)benzene (5.00 g, 21.6 mmol) in Et₂O (50 mL). The mixture was stirred for 30 min at -78 °C and MgBr₂·Et₂O (5.87 g, 22.7 mmol) was added. 20 The mixture was allowed to warm up to 0 °C over 15 min and CuI (420 mg, 2.16 mmol) was added. The mixture was stirred at 0 °C for 5 min and allyl bromide (1.92 mL, 22.7 mmol) was added. The mixture was stirred overnight while warming up to rt. Aq. 1M HCl (1 mL) was added and the mixture was diluted with Et₂O, and washed with aq. 1M HCl (1x). The org. Extracts were dried over 25 MgSO₄, filtered, and the solvents were removed under reduced pressure. The residue was dissolved in acetone (20 mL) and water (10 mL), and TosOH (cat. amount) was added. The mixture was stirred for 5 h at rt, and the solvents were partially removed under reduced pressure. The residue was diluted with Et₂O, and washed with aq. 1M HCl (1x), and aq. sat. NaHCO₃ (1x). The org. extracts were 30 dried over MgSO₄, filtered, and the solvents were removed under reduced pressure. Purification of the residue by FC (Et₂O/petroleum ether 1:49 → 24:1)

yielded 2-allylbenzaldehyde (1.06 g, 34%). This compound was transformed into the title compound following typical procedure J with cyclopropylamine.

Cyclopropyl(2-fluorobenzyl)amine

5

Synthesized according to typical procedure J from 2-fluorobenzaldehyde and cyclopropylamine.

Cyclopropyl-(2-methylbenzyl)amine

10

Synthesized according to typical procedure J from 2-methylbenzaldehyde and cyclopropylamine.

Cyclopropyl-[2-(4-methoxyphenoxy)ethyl]amine

15

Synthesized according to typical procedures C and D from (4-methoxyphenoxy)-acetic acid and cyclopropylamine.

Cyclopropyl-[2-(3-methoxyphenoxy)ethyl]amine

20

Synthesized according to typical procedures C and D from (3-methoxyphenoxy)-acetic acid and cyclopropylamine.

Cyclopropyl-(2-*o*-tolyloxyethyl)amine

25

Synthesized according to typical procedures C and D from *o*-tolyloxyacetic acid and cyclopropylamine.

Cyclopropyl-[2-(3,4-dimethylphenoxy)ethyl]amine

30

Synthesized according to typical procedures C and D from (3,4-dimethylphenoxy)acetic acid and cyclopropylamine.

[2-(2-Chlorophenyl)ethyl]cyclopropylamine

Synthesized according to typical procedures C and D from (2-chlorophenyl)-
5 acetic acid and cyclopropylamine.

Cyclopropyl-[2-(2,3-difluorophenyl)ethyl]amine

Synthesized according to typical procedures C and D from (2,3-difluorophenyl)-
10 acetic acid and cyclopropylamine.

Cyclopropyl-[2-(4-fluorophenyl)ethyl]amine

Synthesized according to typical procedures C and D from (4-fluorophenyl)acetic
15 acid and cyclopropylamine.

Cyclopropyl-(2-*o*-tolylethyl)amine

Synthesized according to typical procedures C and D from *o*-tolylacetic acid and
20 cyclopropylamine.

Cyclopropyl-(2-*p*-tolylethyl)amine

Synthesized according to typical procedures C and D from *p*-tolylacetic acid and
25 cyclopropylamine.

Cyclopropyl-(3,5-dimethoxybenzyl)amine

Synthesized according to typical procedure J from 2,5-dimethoxybenzaldehyde
30 and cyclopropylamine.

(2-Chlorobenzyl)methylamine

See Kihara, M; et al.; *Heterocycles*, 1989, 29, 957.

(2-Chlorobenzyl)isopropylamine

5

Synthesized according to typical procedure J from 2-chlorobenzaldehyde and isopropylamine.

Cyclopropyl-(2-fluoro-5-methoxybenzyl)amine

10

Synthesized according to typical procedure J from 2-fluoro-5-methoxybenzaldehyde and cyclopropylamine.

Cyclopropyl-(3-methoxybenzyl)amine

15

Synthesized according to typical procedure J from 3-methoxybenzaldehyde and cyclopropylamine.

Cyclopropyl-(3,4-dimethoxybenzyl)amine

20

Synthesized according to typical procedure J from 3,4-dimethoxybenzaldehyde and cyclopropylamine.

(2-Chloro-3-trifluoromethylbenzyl)cyclopropylamine

25

Synthesized according to typical procedure J from 2-chloro-3-trifluoromethylbenzaldehyde and cyclopropylamine.

(6-Chlorobenzo[1,3]dioxol-5-ylmethyl)cyclopropylamine

30

Synthesized according to typical procedure J from 6-chlorobenzo[1,3]dioxole-5-carbaldehyde and cyclopropylamine.

Cyclopropyl-(5-fluoro-2-methoxybenzyl)amine

Synthesized according to typical procedure J from 5-fluoro-2-methoxybenzaldehyde and cyclopropylamine.

(2-Chloro-6-fluorobenzyl)cyclopropylamine

Synthesized according to typical procedure J from 2-chloro-6-fluorobenzaldehyde and cyclopropylamine.

(2-Bromobenzyl)cyclopropylamine

Synthesized according to typical procedure J from 2-bromobenzaldehyde and cyclopropylamine.

Cyclopropyl-(2,6-difluorobenzyl)amine

Synthesized according to typical procedure J from 2,6-difluorobenzaldehyde and cyclopropylamine.

Cyclopropyl-(2,3-dimethylbenzyl)amine

Synthesized according to typical procedure J from 2,3-dimethylbenzaldehyde and cyclopropylamine.

Cyclopropyl-(3-fluoro-2-methylbenzyl)amine

Synthesized according to typical procedure J from 3-fluoro-2-methylbenzaldehyde and cyclopropylamine.

Cyclopropyl-(3,5-difluorobenzyl)amine

Synthesized according to typical procedure J from 3,5-difluorobenzaldehyde and cyclopropylamine.

5 **(2-Chloro-3,6-difluorobenzyl)cyclopropylamine**

Synthesized according to typical procedure J from 2-chloro-3,6-difluorobenzaldehyde and cyclopropylamine.

10 **(2,3-Dichlorobenzyl)cyclopropylamine**

Synthesized according to typical procedure J from 2,3-dichlorobenzaldehyde and cyclopropylamine.

15 **Cyclopropyl-(3-trifluoromethoxybenzyl)amine**

Synthesized according to typical procedure J from 3-trifluoromethoxybenzaldehyde and cyclopropylamine.

20 **Cyclopropyl-(3-methylbenzyl)amine**

Synthesized according to typical procedure J from 3-methylbenzaldehyde and cyclopropylamine.

25 **Cyclopropyl-(2,3-difluorobenzyl)amine**

Synthesized according to typical procedure J from 2,3-difluorobenzaldehyde and cyclopropylamine.

30 **(3-Chlorobenzyl)cyclopropylamine**

Synthesized according to typical procedure J from 3-chlorobenzaldehyde and cyclopropylamine.

Cyclopropyl-(4-fluorobenzyl)amine

5

Synthesized according to typical procedure J from 4-fluorobenzaldehyde and cyclopropylamine.

Preparation of other reagents

10

4-Carbamoylbutyric acid

See Melnyk, O., et al.; *J. Org. Chem.*, 2001, 66, 4153.

15 **meso-3,4-Dihydroxytartaric acid anhydride**

A mixture of *meso*-3,4-Dihydroxytartaric acid (1.00 g, 6.67 mmol) and trifluoroacetic acid anhydride (5 mL) was stirred for 2 h at rt. The solvents were removed under reduced pressure and the residue was used as crude product
20 without further purification.

Succinamic acid

See Bellier, B., et al.; *J. Med. Chem.*, 2000, 43, 3614.

25

Specific examples

Example 1

30 (*rac.*)-(2-Methoxyphenyl)acetic acid (*IR**, *5S**)-3-[2-(4-chlorophenyl)acetyl]-
7-{4-[3-(2-methoxybenzyloxy)propoxy]phenyl}-3,9-diazabicyclo[3.3.1]non-6-
en-6-yl methyl ester trifluoroacetate salt

Synthesized according to typical procedure A from bicyclononene N and 4-chlorophenylacetyl chloride, then according to typical procedure E. LC-MS: R_t= 4.57; ES+: 725.35.

5

Example 2

(*rac.*)-(2-Methoxyphenyl)acetic acid (*1R*^{*,} *5S*^{*})-7-{4-[3-(2-methoxybenzyl-oxy)propoxy]phenyl}-3-(quinoxaline-2-carbonyl)-3,9-diazabicyclo[3.3.1]non-10 6-en-6-yl methyl ester trifluoroacetate salt

Synthesized according to typical procedure A from bicyclononene N and 2-quinoxaloyl chloride, then according to typical procedure E. LC-MS: R_t = 4.55; ES+: 728.94.

15

Example 3

(*rac.*)-(2-Methoxyphenyl)acetic acid (*1R*^{*,} *5S*^{*})-7-{4-[3-(2-methoxybenzyl-oxy)propoxy]phenyl}-3-phenylmethanesulfonyl-3,9-diazabicyclo[3.3.1]-non-20 6-en-6-yl methyl ester trifluoroacetate salt

Synthesized according to typical procedure A from bicyclononene N and phenylmethanesulfonyl chloride, then according to typical procedure E. LC-MS: R_t = 4.75; ES+: 727.74.

25

Example 4

(*rac.*)-(2-Methoxyphenyl)acetic acid (*1R*^{*,} *5S*^{*})-7-{4-[3-(2-methoxybenzyl-oxy)propoxy]phenyl}-3-(2-thiophen-2-ylacetyl)-3,9-diazabicyclo[3.3.1]non-30 6-en-6-yl methyl ester trifluoroacetate salt

Synthesized according to typical procedure A from bicyclononene N and thiophen-2-ylacetyl chloride, then according to typical procedure E. LC-MS: R_t = 4.52; ES+: 696.91.

5 **Example 5**

(*rac.*)-(2-Methoxyphenyl)acetic acid (*1R*^{*,} *5S*^{*)}-3-(benzo[*b*]thiophene-3-carbonyl)-7-{4-[3-(2-methoxybenzyloxy)propoxy]phenyl}-3,9-diazabicyclo[3.3.1]non-6-en-6-yl methyl ester trifluoroacetate salt

10

Synthesized according to typical procedure A from bicyclononene N and benzothiophene-3-carbonyl chloride, then according to typical procedure E. LC-MS: R_t = 4.76; ES+: 733.80.

15 **Example 6**

(*rac.*)-(*1R*^{*,} *5S*^{*)}-7-{4-[3-(2-Methoxybenzyloxy)propoxy]phenyl}-6-[2-(2-methoxyphenyl)acetoxymethyl]-3,9-diazabicyclo[3.3.1]non-6-ene-9-carboxylic acid 2,2,2-trichloro-1,1-dimethylethyl ester trifluoroacetate salt

20

Prepared as compound N, but then purified by HPLC. LC-MS: R_t = 5.30; ES+: 774.97.

Example 7

25

(*rac.*)-(*1R*^{*,} *5S*^{*)}-3-Acetyl-7-{4-[3-(2-methoxybenzyloxy)propoxy]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid [2-(2-chlorophenyl)ethyl]-methylamide trifluoroacetate salt

30

Synthesized according to typical procedures H and E from bicyclononene T1 and [2-(2-chlorophenyl)ethyl]methylamine (Jaques B.; Wallace R. G., *Tetrahedron*, 1977, 33, 581). LC-MS: R_t = 0.89; ES+: 632.40.

Example 8

(*rac.*)-(2-Methoxyphenyl)acetic acid (*IR*^{*}, *5S*^{*})-3-acetyl-7-{4-[3-(2-methoxybenzyloxy)propoxy]phenyl}-3,9-diazabicyclo[3.3.1]non-6-en-6-ylmethyl ester trifluoroacetate salt

Synthesized according to typical procedures A and E from bicyclononene N and acetyl chloride. LC-MS: R_t = 0.93; ES+: 615.29.

10

Example 9

(*rac.*)-(*IR*^{*}, *5S*^{*})-3-Acetyl-7-{4-[3-(2-methoxybenzyloxy)propoxy]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid [2-(4-methoxyphenyl)ethyl]-methylamide trifluoroacetate salt

Synthesized according to typical procedures H and E from bicyclononene T1 and [2-(4-methoxyphenyl)ethyl]methylamine (Ho C. Y.; Kukla M. J., *Tetrahedron Lett.* 1997, 38, 2799). LC-MS: R_t = 0.83; ES+: 628.44.

20

Example 10

(*rac.*)-(2-Methoxyphenyl)acetic acid (*IR*^{*}, *5S*^{*})-3-[2-(4-chlorophenyl)acetyl]-7-{6-[3-(2-methoxybenzyloxy)propoxy]pyridin-3-yl}-3,9-diazabicyclo[3.3.1]-non-6-en-6-ylmethyl ester trifluoroacetate salt

Synthesized according to typical procedures A and E from bicyclononene P and 4-chlorophenylacetyl chloride. LC-MS: R_t = 1.03; ES+: 726.44.

30 **Example 11**

(*rac.*)-(IR*, 5S*)-3-Acetyl-7-{4-[3-(2-methoxybenzyloxy)propoxy]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid [2-(4-chlorophenyl)ethyl]-methylamide trifluoroacetate salt

5 Synthesized according to typical procedures H and E from bicyclononene T1 and [2-(4-chlorophenyl)ethyl]methylamine (You Q., et al., *Chem. Res. Chin. Univ.*, 1992, 8, 468). LC-MS: R_t = 0.90; ES+: 632.40.

Example 12

10 **(*rac.*)-(IR*, 5S*)-3-Acetyl-7-{4-[3-(2-methoxybenzyloxy)propoxy]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid [2-(3-chlorophenyl)ethyl]-methylamide trifluoroacetate salt**

15 Synthesized according to typical procedures H and E from bicyclononene T1 and [2-(3-chlorophenyl)ethyl]methylamine. LC-MS: R_t = 0.92; ES+: 632.37.

Example 13

20 **(*rac.*)-(IR*, 5S*)-3-Acetyl-7-{4-[3-(2-methoxybenzyloxy)propoxy]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid ethylphenethylamide trifluoroacetate salt**

25 Synthesized according to typical procedures H and E from bicyclononene T1 and ethylphenethylamine (Cossy J., Rakotoarisoa, H., *Tetrahedron Lett.*, 2000, 41, 2097). LC-MS: R_t = 0.89; ES+: 612.46.

Example 14

30 **(*rac.*)-(IR*, 5S*)-3-Acetyl-7-{4-[3-(2-methoxybenzyloxy)propoxy]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid [2-(3-methoxyphenyl)ethyl]-methylamide trifluoroacetate salt**

Synthesized according to typical procedures H and E from bicyclononene T1 and [2-(3-methoxyphenyl)ethyl]methylamine. LC-MS: R_t = 0.88; ES+: 628.41.

5 **Example 15**

(*rac.*)-(2-Methoxyphenyl)acetic acid (*1R*^{*}, *5S*^{*})-7-{4-[3-(2-methoxybenzyl
oxy)propoxy]phenyl}-3-methyl-3,9-diazabicyclo[3.3.1]non-6-en-6-ylmethyl
ester trifluoroacetate salt

10

A sol. of bicyclononene N (0.05 mmol) in CH₂Cl₂ (2 mL) was cooled to 0 °C. DIPEA (0.10 mmol) and methyl iodide (0.10 mmol) were added. The mixture was stirred at 0 °C for 2 h, then overnight at rt. Methyl iodide (0.50 mmol) and DIPEA (0.15 mmol) were added again and stirring was continued for 4 h at rt. EtOAc was added and the mixture was washed with water. The org. extracts were separated, dried over MgSO₄ and filtered. The solvents were removed under reduced pressure and the residue was processed further according to general procedure E. LC-MS: R_t = 4.04; ES+: 587.38.

20 **Example 16**

(*rac.*)-(1*R*^{*}, 5*S*^{*})-3-Acetyl-7-{4-[3-(2-methoxybenzyl)oxy]propoxy}phenyl}-3,9-
diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid [2-(3,4-dimethoxyphenyl)-
ethyl]methylamide trifluoroacetate salt

25

Synthesized according to typical procedures H and E from bicyclononene T1 and [2-(3,4-dimethoxyphenyl)ethyl]methylamine. LC-MS: R_t = 0.89; ES+: 644.48.

Example 17

30

(*rac.*)-(2-Methoxyphenyl)acetic acid (*1R*^{*}, *5S*^{*})-7-{4-[3-(2-methoxybenzyl-oxy)propoxy]phenyl}-3,9-diazabicyclo[3.3.1]non-6-en-6-ylmethyl ester trifluoroacetate salt

5 Synthesized according to typical procedure E from bicyclononene N. LC-MS: R_t = 0.83; ES+: 573.29.

Example 18

10 **(*rac.*)-N-((*1R*^{*}, *5S*^{*})-3-Acetyl-7-{4-[3-(2-methoxybenzyloxy)propoxy]phenyl}-3,9-diazabicyclo[3.3.1]non-6-en-6-ylmethyl)-2-(2-methoxyphenyl)-*N*-methyl-acetamide trifluoroacetate salt**

15 Synthesized according to typical procedures G and E from bicyclononene Z and (2-methoxyphenyl)acetic acid. LC-MS: R_t = 3.98; ES+: 628.63.

Example 19

20 **(*rac.*)-(*1R*^{*}, *5S*^{*})-3-Acetyl-7-{4-[3-(2-methoxybenzyloxy)propoxy]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid methyl(3-phenylpropyl)amide trifluoroacetate salt**

25 Synthesized according to typical procedures H and E from bicyclononene T1 and *N*-methyl(3-phenylpropyl)amine (Lavastre I.; *et al.*, *Bull. Soc. Chim. Fr.*, 1995, 132, 188). LC-MS: R_t = 0.88; ES+: 612.45.

Example 20

30 **(*rac.*)-(*1R*^{*}, *5S*^{*})-3-Acetyl-7-{4-[3-(2-methoxybenzyloxy)propoxy]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid [2-(2-methoxyphenyl)ethyl]-methylamide trifluoroacetate**

Synthesized according to typical procedures H and E from bicyclononene **T1** and [2-(2-methoxyphenyl)ethyl]methylamine. LC-MS: R_t = 0.88; ES+: 628.45.

Example 21

5

(*rac.*)-(IR*, 5S*)-3-Acetyl-7-{4-[3-(2-methoxybenzyloxy)propoxy]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid benzylmethyleamide trifluoroacetate

10 Synthesized according to typical procedures H and E from bicyclononene **T1** and N-methylbenzylamine. LC-MS: R_t = 0.84; ES+: 684.41.

Example 22

15

(*rac.*)-(2-Methoxyphenyl)acetic acid (IR*, 5S*)-3-acetyl-7-{4-[3-(2-chlorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-en-6-ylmethyl ester trifluoroacetate

20

Synthesized according to typical procedures F and E from bicyclononene **R2** and 2-chlorophenol. LC-MS: R_t = 0.89; ES+: 589.34.

Example 23

25

(*rac.*)-(2-Methoxyphenyl)acetic acid (IR*, 5S*)-3-acetyl-7-{6-[3-(2-methoxybenzyloxy)propoxy]pyridin-3-yl}-3,9-diazabicyclo[3.3.1]non-6-en-6-ylmethyl ester trifluoroacetate salt

30

Example 24

Synthesized according to typical procedures A and E from bicyclononene **P** and acetyl chloride. LC-MS: R_t = 0.90; ES+: 616.44.

(rac.)-(2-Methoxyphenyl)acetic acid ($IR^, 5S^*$)-3-acetyl-7-{4-[3-(2-bromo-
phenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-en-6-ylmethyl ester
trifluoroacetate salt*

- 5 Synthesized according to typical procedures F and E from bicyclononene R2 and
2-bromophenol. LC-MS: $R_t = 0.92$; ES+: 633.26.

Example 25

- 10 *(rac.)-(2-Methoxyphenyl)acetic acid ($IR^*, 5S^*$)-3-[2-(4-chlorophenyl)ethyl]-7-
{4-[3-(2-methoxybenzyloxy)propanoyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-en-
6-yl- methyl ester trifluoroacetate salt*

To a sol. of bicyclononene N (0.05 mmol) in CH_2Cl_2 (2 mL) NaBH_3OAc (0.065
15 mmol) and (4-chlorophenyl)acetaldehyde (Zhuangyu Z., et al., *Synthesis*, 1991,
539, 0.065 mmol) were added. The mixture was stirred overnight. The solvent
was removed under reduced pressure and the residue treated according to typical
procedure E. LC-MS: $R_t = 4.88$; ES+: 711.47.

20 **Example 26**

(rac.)-(2-Methoxyphenyl)acetic acid ($IR^, 5S^*$)-3-acetyl-7-{4-[2-(3-chloro-
phenoxy)ethyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-en-6-ylmethyl ester
trifluoroacetate salt*

- 25 Synthesized according to typical procedures F and E from bicyclononene R1 and
3-chlorophenol. LC-MS: $R_t = 4.17$; ES+: 575.62.

Example 27

(*rac.*)-(2-Methoxyphenyl)acetic acid (*1R*^{*,}, *5S*^{*})-3-acetyl-7-{4-[3-(3-chlorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-en-6-ylmethyl ester trifluoro-acetate salt

- 5 Synthesized according to typical procedures F and E from bicyclononene R2 and 3-chlorophenol. LC-MS: R_t = 0.90; ES+: 589.33.

Example 28

- 10 (*rac.*)-(2-Methoxyphenyl)acetic acid (*1R*^{*,}, *5S*^{*})-3-acetyl-7-{4-[2-(2-chlorophenoxy)ethyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-en-6-ylmethyl ester trifluoroacetate salt

- 15 Synthesized according to typical procedures F and E from bicyclononene R1 and 2-chlorophenol. LC-MS: R_t = 0.94; ES+: 575.34.

Example 29

- 20 (*rac.*)-(*1R*^{*,} *5S*^{*})-3-Acetyl-7-{4-[3-(2-methoxybenzyloxy)propoxy]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid phenethylamide trifluoroacetate salt

- 25 Synthesized according to typical procedures H and E from bicyclononene T1 and phenethylamine. LC-MS: R_t = 0.91; ES+: 584.44.

25

Example 30

- 30 (*rac.*)-(2-Methoxyphenyl)acetic acid (*1R*^{*,}, *5S*^{*})-3-acetyl-7-[4-(3-phenoxy-propyl)phenyl]-3,9-diazabicyclo[3.3.1]non-6-en-6-ylmethyl ester trifluoro-acetate salt

Synthesized according to typical procedures F and E from bicyclononene R2 and phenol. LC-MS: R_t = 0.87; ES+: 555.31.

Example 31

5

1:1-Mixture of (2R)- and (2S)-N-((1R*, 5S*)-3-acetyl-7-{4-[3-(2-methoxybenzyloxy)propoxy]phenyl}-3,9-diazabicyclo[3.3.1]non-6-en-6-ylmethyl)-N-methyl-2-phenylpropionamide trifluoroacetate salt

10 Synthesized according to typical procedures G and E from bicyclononene Z and (rac.)-2-phenylpropionic acid. LC-MS: R_t = 0.87; ES+: 612.42.

Example 32

15 **(rac.)-(1R*, 5S*)-3-Acetyl-7-{4-[3-(2-methoxybenzyloxy)propoxy]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid [2-(2-methoxyphenyl)ethyl]-amide trifluoroacetate salt**

20 Synthesized according to typical procedures H and E from bicyclononene T1 and 2-(2-methoxyphenyl)ethylamine. LC-MS: R_t = 4.06; ES+: 614.35.

Example 33

25 **(rac.)-(1R*, 5S*)-3-Acetyl-7-{4-[2-(2-chlorophenoxy)ethyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid methylphenethylamide formate salt**

Synthesized according to typical procedures F and E from bicyclononene U2 and 2-chlorophenol. LC-MS: R_t = 0.84; ES+: 558.24.

30 **Example 34**

(*rac.*)-(1*R*^{*}, 5*S*^{*})-3-Acetyl-7-{4-[2-(2-ethoxyphenoxy)ethyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid methylphenethylamide formate salt

Synthesized according to typical procedures F and E from bicyclononene U2 and
5 2-ethoxyphenol. LC-MS: R_t = 0.84; ES+: 568.30.

Example 35

(*rac.*)-(1*R*^{*}, 5*S*^{*})-3-Acetyl-7-{4-[2-(2-acetylphenoxy)ethyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid methylphenethylamide formate salt

Synthesized according to typical procedures F and E from bicyclononene U2 and
2-acetylphenol. LC-MS: R_t = 1.13; ES+: 566.29.

15 Example 36

(*rac.*)-(1*R*^{*}, 5*S*^{*})-3-Acetyl-7-{4-[2-*o*-toluoxyethyl]phenyl}-3,9-diazabicyclo[3.3.1]-non-6-ene-6-carboxylic acid methylphenethylamide formate salt

20 Synthesized according to typical procedures F and E from bicyclononene U2 and
2-methylphenol. LC-MS: R_t = 1.18; ES+: 538.27.

Example 37

25 (*rac.*)-(1*R*^{*}, 5*S*^{*})-3-Acetyl-7-{4-[2-(3-methoxyphenoxy)ethyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid methylphenethylamide formate salt

Synthesized according to typical procedures F and E from bicyclononene U2 and
30 3-methoxyphenol. LC-MS: R_t = 1.15; ES+: 554.28.

Example 38

(rac.)-(1R, 5S*)-3-Acetyl-7-{4-[2-(3-trifluoromethoxyphenoxy)ethyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid methylphenethylamide formate salt*

5

Synthesized according to typical procedures F and E from bicyclononene U2 and 3-trifluoromethoxyphenol. LC-MS: R_t = 1.19; ES+: 608.28.

Example 39

10

(rac.)-(1R, 5S*)-3-Acetyl-7-{4-[2-m-toluoxoxyethyl]phenyl}-3,9-diazabicyclo[3.3.1]-non-6-ene-6-carboxylic acid methylphenethylamide formate salt*

15

Synthesized according to typical procedures F and E from bicyclononene U2 and 3-methylphenol. LC-MS: R_t = 1.17; ES+: 538.26.

Example 40

20

(rac.)-(1R, 5S*)-3-Acetyl-7-{4-[2-(3-isopropylphenoxy)ethyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid methylphenethylamide formate salt*

25

Synthesized according to typical procedures F and E from bicyclononene U2 and 3-isopropylphenol. LC-MS: R_t = 1.19; ES+: 608.28.

Example 41

30

(rac.)-(1R, 5S*)-3-Acetyl-7-{4-[3-(2-chlorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid methylphenethylamide trifluoroacetate salt*

Synthesized according to typical procedures F and E from bicyclononene U3 and 2-chlorophenol. LC-MS: R_t = 0.87; ES+: 572.13.

Example 42

5

(*rac.*)-(1*R*^{*}, 5*S*^{*})-3-Acetyl-7-{4-[3-(2-bromophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid methylphenethylamide formate salt

10 Synthesized according to typical procedures F and E from bicyclononene U3 and 2-bromophenol. LC-MS: R_t = 0.79; ES+: 616.11.

Example 43

15

(*rac.*)-(1*R*^{*}, 5*S*^{*})-3-Acetyl-7-{4-[3-(2-fluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid methylphenethylamide formate salt

Synthesized according to typical procedures F and E from bicyclononene U3 and 2-fluorophenol. LC-MS: R_t = 0.75; ES+: 556.20.

20

Example 44

25

(*rac.*)-(1*R*^{*}, 5*S*^{*})-3-Acetyl-7-{4-[3-(2-acetylphenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid methylphenethylamide formate salt

25

Synthesized according to typical procedures F and E from bicyclononene U3 and 2-acetylphenol. LC-MS: R_t = 0.71; ES+: 580.20.

Example 45

30

(rac.)-(IR, 5S*)-3-Acetyl-7-{4-[3-o-toluoxypropyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid methylphenethylamide trifluoroacetate salt*

- 5 Synthesized according to typical procedures F and E from bicyclononene U3 and 2-methylphenol. LC-MS: R_t = 0.90; ES+: 552.20.

Example 46

- 10 *(rac.)-(IR*, 5S*)-3-Acetyl-7-{4-[3-(3-methoxyphenoxy)propyl]phenyl}-3,9-di-azabicyclo[3.3.1]non-6-ene-6-carboxylic acid methylphenethylamide trifluoroacetate salt*

Synthesized according to typical procedures F and E from bicyclononene U3 and 15 3-methoxyphenol. LC-MS: R_t = 0.86; ES+: 568.21.

Example 47

- 20 *(rac.)-(IR*, 5S*)-3-Acetyl-7-{4-[3-m-toluoxypropyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid methylphenethylamide formate salt*

Synthesized according to typical procedures F and E from bicyclononene U3 and 3-methylphenol. LC-MS: R_t = 0.79; ES+: 552.21.

25 **Example 48**

(rac.)-(IR, 5S*)-3-Acetyl-7-{4-[3-(3-chlorophenoxy)propyl]phenyl}-3,9-di-azabicyclo[3.3.1]non-6-ene-6-carboxylic acid methylphenethylamide trifluoroacetate salt*

- 30 Synthesized according to typical procedures F and E from bicyclononene U3 and 3-chlorophenol. LC-MS: R_t = 4.44; ES+: 572.32.

Example 49

5 **(*rac.*)-(1*R*^{*}, 5*S*^{*})-3-Acetyl-7-{4-[2-(3-bromophenoxy)ethyl]phenyl}-3,9-diaza-**
bicyclo[3.3.1]non-6-ene-6-carboxylic acid methylphenethylamide formate salt

Synthesized according to typical procedures F and E from bicyclononene U2 and 3-bromophenol. LC-MS: R_t = 0.78; ES+: 602.10.

10 **Example 50**

15 **(*rac.*)-(1*R*^{*}, 5*S*^{*})-3-Acetyl-7-{4-[2-(2,3-dichlorophenoxy)ethyl]phenyl}-3,9-di-**
azabicyclo[3.3.1]non-6-ene-6-carboxylic acid methylphenethylamide formate
salt

15 Synthesized according to typical procedures F and E from bicyclononene U2 and 2,3-dichlorophenol. LC-MS: R_t = 0.78; ES+: 592.10.

20 **Example 51**

20 **(*rac.*)-(1*R*^{*}, 5*S*^{*})-3-Acetyl-7-{4-[2-(2-chloro-3-trifluoromethylphenoxy)ethyl]-**
phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid methylphenethyl-
amide formate salt

25 Synthesized according to typical procedures F and E from bicyclononene U2 and 2-chloro-3-trifluoromethylphenol. LC-MS: R_t = 0.80; ES+: 626.11.

25 **Example 52**

30 **(*rac.*)-(1*R*^{*}, 5*S*^{*})-3-Acetyl-7-{4-[2-(2,3-difluorophenoxy)ethyl]phenyl}-3,9-di-**
azabicyclo[3.3.1]non-6-ene-6-carboxylic acid methylphenethylamide formate
salt

Synthesized according to typical procedures F and E from bicyclononene **U2** and 2,3-difluorophenol. LC-MS: $R_t = 0.74$; ES+: 560.13.

5 **Example 53**

(*rac.*)-(IR*, 5S*)-3-Acetyl-7-{4-[2-(2,3-dimethylphenoxy)ethyl]phenyl}-3,9-di-azabicyclo[3.3.1]non-6-ene-6-carboxylic acid methylphenethylamide formate salt

10

Synthesized according to typical procedures F and E from bicyclononene **U2** and 2,3-dimethylphenol. LC-MS: $R_t = 0.79$; ES+: 559.19.

Example 54

15

(*rac.*)-(IR*, 5S*)-3-Acetyl-7-{4-[2-(2-ethylphenoxy)ethyl]phenyl}-3,9-diaza-bicyclo[3.3.1]non-6-ene-6-carboxylic acid methylphenethylamide formate salt

20

Synthesized according to typical procedures F and E from bicyclononene **U2** and 2-ethylphenol. LC-MS: $R_t = 0.80$; ES+: 552.19.

Example 55

25

(*rac.*)-(IR*, 5S*)-3-Acetyl-7-{4-[3-(3-bromophenoxy)propyl]phenyl}-3,9-di-azabicyclo[3.3.1]non-6-ene-6-carboxylic acid methylphenethylamide trifluoroacetate salt

Synthesized according to typical procedures F and E from bicyclononene **U3** and 3-bromophenol. LC-MS: $R_t = 0.92$; ES+: 616.05.

30

Example 56

(*rac.*)-(1*R*^{*}, 5*S*^{*})-3-Acetyl-7-{4-[3-(2,3-dichlorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid methylphenethylamide trifluoroacetate salt

- 5 Synthesized according to typical procedures F and E from bicyclononene U3 and 2,3-dichlorophenol. LC-MS: R_t = 0.93; ES+: 606.12.

Example 57

10 **(*rac.*)-(1*R*^{*}, 5*S*^{*})-3-Acetyl-7-{4-[3-(2-chloro-3-trifluoromethylphenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid methylphenethylamide trifluoroacetate salt**

- 15 Synthesized according to typical procedures F and E from bicyclononene U3 and 2-chloro-3-trifluoromethylphenol. LC-MS: R_t = 0.93; ES+: 640.10.

Example 58

20 **(*rac.*)-(1*R*^{*}, 5*S*^{*})-3-Acetyl-7-{4-[3-(2,3-difluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid methylphenethylamide trifluoroacetate salt**

- 25 Synthesized according to typical procedures F and E from bicyclononene U3 and 2,3-difluorophenol. LC-MS: R_t = 0.88; ES+: 574.17.

25

Example 59

30 **(*rac.*)-(1*R*^{*}, 5*S*^{*})-3-Acetyl-7-{4-[3-(2,3-dimethylphenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid methylphenethylamide trifluoroacetate salt**

Synthesized according to typical procedures F and E from bicyclononene U3 and 2,3-dimethylphenol. LC-MS: R_t = 0.93; ES+: 566.22.

Example 60

5

(*rac.*)-(IR*, 5S*)-3-Acetyl-7-{4-[3-(2-ethylphenoxy)propyl]phenyl}-3,9-diaza-bicyclo[3.3.1]non-6-ene-6-carboxylic acid methylphenethylamide trifluoro-acetate salt

10 Synthesized according to typical procedures F and E from bicyclononene U3 and 2-ethylphenol. LC-MS: R_t = 0.93; ES+: 566.22.

Example 61

15

(*rac.*)-(IR*, 5S*)-3-Acetyl-7-{4-[3-(2-isopropylphenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid methylphenethylamide formate salt

20

Synthesized according to typical procedures F and E from bicyclononene U3 and 2-isopropylphenol. LC-MS: R_t = 0.86; ES+: 580.22.

Example 62

25

(*rac.*)-(IR*, 5S*)-3-Acetyl-7-{4-[3-(2-*tert*-butylphenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid methylphenethylamide formate salt

30

Synthesized according to typical procedures F and E from bicyclononene U3 and 2-*tert*-butylphenol. LC-MS: R_t = 0.89; ES+: 594.24.

Example 63

(rac.)-(1R, 5S*)-3-Acetyl-7-{4-[3-(4-chloro-2-methoxyphenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid methylphenethylamide formate salt*

5 Synthesized according to typical procedures F and E from bicyclononene U3 and 4-chloro-2-methoxyphenol. LC-MS: R_t = 0.77; ES+: 602.14.

Example 64

10 *(rac.)-(1R*, 5S*)-3-Acetyl-7-{4-[3-(2,4-dichlorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid methylphenethylamide formate salt*

15 Synthesized according to typical procedures F and E from bicyclononene U3 and 2,4-dichlorophenol. LC-MS: R_t = 0.84; ES+: 606.08.

Example 65

20 *(rac.)-(1R*, 5S*)-3-Acetyl-7-{4-[3-(4-fluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid methylphenethylamide formate salt*

Synthesized according to typical procedures F and E from bicyclononene U3 and 4-fluorophenol. LC-MS: R_t = 0.76; ES+: 556.19.

25 **Example 66**

(rac.)-(1R, 5S*)-3-Acetyl-7-{4-[3-(2-*tert*-butyl-4-methylphenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid methylphenethylamide formate salt*

30 Synthesized according to typical procedures F and E from bicyclononene U3 and 2-*tert*-butyl-4-methylphenol. LC-MS: R_t = 0.93; ES+: 608.25.

Example 67

5 (rac.)-(*1R*^{*,} *5S*^{*)}-3-Acetyl-7-{4-[3-(2-bromo-5-fluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid methylphenethylamide formate salt

Synthesized according to typical procedures F and E from bicyclononene U3 and 2-bromo-5-fluorophenol. LC-MS: R_t = 0.80; ES+: 634.05.

10

Example 68

15 (rac.)-(*1R*^{*,} *5S*^{*)}-3-Acetyl-7-{4-[3-(2,5-difluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid methylphenethylamide formate salt

Synthesized according to typical procedures F and E from bicyclononene U3 and 2,5-difluorophenol. LC-MS: R_t = 0.76; ES+: 574.14.

20 **Example 69**

(rac.)-(*1R*^{*,} *5S*^{*)}-3-Acetyl-7-{4-[3-(2-chloro-5-methylphenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid methylphenethylamide formate salt

25

Synthesized according to typical procedures F and E from bicyclononene U3 and 2-chloro-5-methylphenol. LC-MS: R_t = 0.82; ES+: 586.16.

Example 70

30

(rac.)-(1R, 5S*)-3-Acetyl-7-{4-[3-(2,5-dimethylphenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid methylphenethylamide formate salt*

- 5 Synthesized according to typical procedures F and E from bicyclononene U3 and
2,5-dimethylphenol. LC-MS: R_t = 0.83; ES+: 566.20.

Example 71

- 10 *(rac.)-(1R*, 5S*)-3-Acetyl-7-{4-[2-(2-isopropylphenoxy)ethyl]phenyl}-3,9-di-
azabicyclo[3.3.1]non-6-ene-6-carboxylic acid methylphenethylamide formate
salt*

15 Synthesized according to typical procedures F and E from bicyclononene U2 and
2-isopropylphenol. LC-MS: R_t = 0.82; ES+: 566.22.

Example 72

- 20 *(rac.)-(1R*, 5S*)-3-Acetyl-7-{4-[2-(2-*tert*-butylphenoxy)ethyl]phenyl}-3,9-di-
azabicyclo[3.3.1]non-6-ene-6-carboxylic acid methylphenethylamide formate
salt*

25 Synthesized according to typical procedures F and E from bicyclononene U2 and
2-*tert*-butylphenol. LC-MS: R_t = 0.85; ES+: 580.26.

Example 73

- 30 *(rac.)-(1R*, 5S*)-3-Acetyl-7-{4-[2-(2-propionylphenoxy)ethyl]phenyl}-3,9-di-
azabicyclo[3.3.1]non-6-ene-6-carboxylic acid methylphenethylamide formate
salt*

Synthesized according to typical procedures F and E from bicyclononene U2 and 2-propionylphenol. LC-MS: R_t = 0.72; ES+: 580.21.

Example 74

5

(rac.)-(IR*, 5S*)-3-Acetyl-7-{4-[2-(2,4-dichlorophenoxy)ethyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid methylphenethylamide formate salt

10

Synthesized according to typical procedures F and E from bicyclononene U2 and 2,4-dichlorophenol. LC-MS: R_t = 0.80; ES+: 592.09.

Example 75

15

(rac.)-(IR*, 5S*)-3-Acetyl-7-{4-[2-(2-*tert*-butyl-4-methylphenoxy)ethyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid methylphenethylamide formate salt

20

Synthesized according to typical procedures F and E from bicyclononene U2 and 2-*tert*-butyl-4-methylphenol. LC-MS: R_t = 0.88; ES+: 594.27.

Example 76

25

(rac.)-(IR*, 5S*)-3-Acetyl-7-{4-[2-(4-methoxyphenoxy)ethyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid methylphenethylamide formate salt

30

Synthesized according to typical procedures F and E from bicyclononene U2 and 4-methoxyphenol. LC-MS: R_t = 0.71; ES+: 554.18.

Example 77

(*rac.*)-(1*R*^{*}, 5*S*^{*})-3-Acetyl-7-{4-[2-(2-bromo-5-fluorophenoxy)ethyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid methylphenethylamide formate salt

5

Synthesized according to typical procedures F and E from bicyclononene U2 and 2-bromo-5-fluorophenol. LC-MS: R_t = 0.76; ES+: 620.09.

Example 78

10

(*rac.*)-(1*R*^{*}, 5*S*^{*})-3-Acetyl-7-{4-[2-(2,5-difluorophenoxy)ethyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid methylphenethylamide formate salt

15

Synthesized according to typical procedures F and E from bicyclononene U2 and 2,5-difluorophenol. LC-MS: R_t = 0.73; ES+: 560.17.

Example 79

20

(*rac.*)-(1*R*^{*}, 5*S*^{*})-3-Acetyl-7-{4-[2-(2-chloro-5-methylphenoxy)ethyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid methylphenethylamide formate salt

25

Synthesized according to typical procedures F and E from bicyclononene U2 and 2-chloro-5-methylphenol. LC-MS: R_t = 0.78; ES+: 572.13.

Example 80

30

(*rac.*)-(1*R*^{*}, 5*S*^{*})-3-Acetyl-7-{4-[2-(2-methoxy-5-methylphenoxy)ethyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid methylphenethylamide formate salt

Synthesized according to typical procedures F and E from bicyclononene **U2** and 2-methoxy-5-methylphenol. LC-MS: R_t = 0.72; ES+: 568.19.

Example 81

5

(*rac.*)-(1*R*^{*}, 5*S*^{*})-3-Acetyl-7-{4-[2-(2,5-dimethylphenoxy)ethyl]phenyl}-3,9-di-azabicyclo[3.3.1]non-6-ene-6-carboxylic acid methylphenethylamide formate salt

10 Synthesized according to typical procedures F and E from bicyclononene **U2** and 2,5-dimethylphenol. LC-MS: R_t = 0.90; ES+: 552.24.

Example 82

15 (*rac.*)-(1*R*^{*}, 5*S*^{*})-3-Acetyl-7-{4-[2-(2-chlorophenoxy)ethoxy]phenyl}-3,9-di-azabicyclo[3.3.1]non-6-ene-6-carboxylic acid methylphenethylamide formate salt

20 Synthesized according to typical procedures F and E from bicyclononene **U1** and 2-(2-chlorophenoxy)ethanol. LC-MS: R_t = 0.75; ES+: 574.15.

Example 83

25 (*rac.*)-(1*R*^{*}, 5*S*^{*})-3-Acetyl-7-{4-[2-(3-methylphenoxy)ethoxy]phenyl}-3,9-di-azabicyclo[3.3.1]non-6-ene-6-carboxylic acid methylphenethylamide formate salt

Synthesized according to typical procedures F and E from bicyclononene **U1** and 2-(3-methylphenoxy)ethanol. LC-MS: R_t = 0.76; ES+: 554.18.

30

Example 84

(*rac.*)-(1*R*^{*}, 5*S*^{*})-3-Acetyl-7-{4-[2-(2-chlorophenyl)ethoxy]phenyl}-3,9-diaza-bicyclo[3.3.1]non-6-ene-6-carboxylic acid methylphenethylamide formate salt

Synthesized according to typical procedures F and E from bicyclononene U1 and
5 2-(chlorophenyl)ethanol. LC-MS: R_t = 0.77; ES+: 558.13.

Example 85

(*rac.*)-(1*R*^{*}, 5*S*^{*})-3-Acetyl-7-{4-[2-(3-chlorophenyl)ethoxy]phenyl}-3,9-diaza-bicyclo[3.3.1]non-6-ene-6-carboxylic acid methylphenethylamide formate salt

Synthesized according to typical procedures F and E from bicyclononene U1 and
2-(3-chlorophenyl)ethanol. LC-MS: R_t = 0.77; ES+: 558.14.

15 Example 86

(*rac.*)-(1*R*^{*}, 5*S*^{*})-3-Acetyl-7-{4-[2-(2-methoxyphenyl)ethoxy]phenyl}-3,9-di-azabicyclo[3.3.1]non-6-ene-6-carboxylic acid methylphenethylamide trifluoroacetate salt

20 Synthesized according to typical procedures F and E from bicyclononene U1 and
2-(2-methoxyphenyl)ethanol. LC-MS: R_t = 0.85; ES+: 554.21.

Example 87

25 (*rac.*)-(1*R*^{*}, 5*S*^{*})-3-Acetyl-7-{4-[2-(2,5-dichlorophenoxy)ethyl]phenyl}-3,9-di-azabicyclo[3.3.1]non-6-ene-6-carboxylic acid methylphenethylamide formate salt

30 Synthesized according to typical procedures F and E from bicyclononene U2 and
2,5-dichlorophenol. LC-MS: R_t = 0.79; ES+: 592.07.

Example 88

(*rac.*)-(1*R*^{*,}, 5*S*^{*})-3-Acetyl-7-{4-[2-(2,3,6-trifluorophenoxy)ethyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid methylphenethylamide
5 formate salt

Synthesized according to typical procedures F and E from bicyclononene **U2** and 2,3,6-trifluorophenol. LC-MS: R_t = 0.74; ES+: 578.14.

10 Example 89

(*rac.*)-(1*R*^{*,}, 5*S*^{*})-3-Acetyl-7-{4-[2-(3,5-dimethylphenoxy)ethyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid methylphenethylamide formate salt

15 Synthesized according to typical procedures F and E from bicyclononene **U2** and 3,5-dimethylphenol. LC-MS: R_t = 0.80; ES+: 552.20.

Example 90

20 (*rac.*)-(1*R*^{*,}, 5*S*^{*})-3-Acetyl-7-{4-[2-(3-chlorophenoxy)ethyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid methylphenethylamide formate salt

25 Synthesized according to typical procedures F and E from bicyclononene **U2** and 3-chlorophenol. LC-MS: R_t = 0.77; ES+: 558.15.

Example 91

30 (*rac.*)-(1*R*^{*,}, 5*S*^{*})-3-Acetyl-7-{4-[2-(3-trifluoromethylphenoxy)ethyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid methylphenethylamide formate salt

Synthesized according to typical procedures F and E from bicyclononene **U2** and 3-trifluoromethylphenol. LC-MS: R_t = 0.78; ES+: 592.17.

5 **Example 92**

(*rac.*)-(1*R*^{*}, 5*S*^{*})-3-Acetyl-7-{4-[2-(4-*tert*-butyl-2-methylphenoxy)ethyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid methylphenethylamide formate salt

10

Synthesized according to typical procedures F and E from bicyclononene **U2** and 4-*tert*-butyl-2-methylphenol. LC-MS: R_t = 0.88; ES+: 594.22.

Example 93

15

(*rac.*)-(1*R*^{*}, 5*S*^{*})-3-Acetyl-7-{4-[2-(3,4-dichlorophenoxy)ethyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid methylphenethylamide formate salt

20

Synthesized according to typical procedures F and E from bicyclononene **U2** and 3,4-dichlorophenol. LC-MS: R_t = 0.81; ES+: 592.12.

Example 94

25

(*rac.*)-(1*R*^{*}, 5*S*^{*})-3-Acetyl-7-{4-[2-(4-bromo-3-methylphenoxy)ethyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid methylphenethylamide formate salt

Synthesized according to typical procedures F and E from bicyclononene **U2** and

30

4-bromo-3-methylphenol. LC-MS: R_t = 0.81; ES+: 616.12.

Example 95

(*rac.*)-(1*R*^{*, 5*S*})-3-Acetyl-7-{4-[2-(3,4-dimethylphenoxy)ethyl]phenyl}-3,9-di-azabicyclo[3.3.1]non-6-ene-6-carboxylic acid methylphenethylamide formate salt

5

Synthesized according to typical procedures F and E from bicyclononene U2 and 3,4-dimethylphenol. LC-MS: R_t = 0.78; ES+: 552.18.

Example 96

10

(*rac.*)-(1*R*^{*, 5*S*})-3-Acetyl-7-{4-[2-(3,5-dichlorophenoxy)ethyl]phenyl}-3,9-di-azabicyclo[3.3.1]non-6-ene-6-carboxylic acid methylphenethylamide formate salt

15

Synthesized according to typical procedures F and E from bicyclononene U2 and 3,5-dichlorophenol. LC-MS: R_t = 0.82; ES+: 592.10.

Example 97

20

(*rac.*)-(1*R*^{*, 5*S*})-3-Acetyl-7-{4-[2-(3,5-dimethylphenoxy)ethyl]phenyl}-3,9-di-azabicyclo[3.3.1]non-6-ene-6-carboxylic acid methylphenethylamide formate salt

25

Synthesized according to typical procedures F and E from bicyclononene U2 and 3,5-dimethylphenol. LC-MS: R_t = 0.79; ES+: 552.20.

Example 98

30

(*rac.*)-(1*R*^{*, 5*S*})-3-Acetyl-7-{4-[2-(3,5-dimethoxyphenoxy)ethyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid methylphenethylamide formate salt

Synthesized according to typical procedures F and E from bicyclononene **U2** and 3,5-dimethoxyphenol. LC-MS: R_t = 0.72; ES+: 584.19.

Example 99

5

(*rac.*)-(IR*, SS*)-3-Acetyl-7-{4-[2-(2-chloro-4,5-dimethylphenoxy)ethyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid methylphenethylamide formate salt

10 Synthesized according to typical procedures F and E from bicyclononene **U2** and 2-dichloro-4,5-dimethylphenol. LC-MS: R_t = 0.81; ES+: 586.17.

Example 100

15 (*rac.*)-(IR*, SS*)-3-Acetyl-7-{4-[2-(2,3,5-trimethylphenoxy)ethyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid methylphenethylamide formate salt

20 Synthesized according to typical procedures F and E from bicyclononene **U2** and 2,3,5-trimethylphenol. LC-MS: R_t = 0.82; ES+: 566.21.

Example 101

25 (*rac.*)-(IR*, SS*)-3-Acetyl-7-{4-[3-(2,5-dichlorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid methylphenethylamide formate salt

Synthesized according to typical procedures F and E from bicyclononene **U3** and 2,5-dichlorophenol. LC-MS: R_t = 0.83; ES+: 606.12.

30

Example 102

(rac.)-(1R, 5S*)-3-Acetyl-7-{4-[3-(2-acetyl-5-fluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid methylphenethylamide formate salt*

- 5 Synthesized according to typical procedures F and E from bicyclononene U3 and 2-acetyl-5-fluorophenol. LC-MS: R_t = 0.73; ES+: 598.18.

Example 103

- 10 *(rac.)-(1R*, 5S*)-3-Acetyl-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid methylphenethylamide trifluoroacetate salt*

Synthesized according to typical procedures F and E from bicyclononene U3 and 15 2,3,6-trifluorophenol. LC-MS: R_t = 0.88; ES+: 592.19.

Example 104

- 20 *(rac.)-(1R*, 5S*)-3-Acetyl-7-{4-[3-(2,4-dimethylphenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid methylphenethylamide formate salt*

Synthesized according to typical procedures F and E from bicyclononene U3 and 25 2,4-dimethylphenol. LC-MS: R_t = 0.83; ES+: 566.22.

25

Example 105

- 30 *(rac.)-(1R*, 5S*)-3-Acetyl-7-{4-[3-(2-*tert*-butyl-6-methylphenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid methylphenethylamide formate salt*

Synthesized according to typical procedures F and E from bicyclononene U3 and 2-*tert*-butyl-6-methylphenol. LC-MS: R_t = 0.88; ES+: 608.27.

Example 106

5

(*rac.*)-(1*R*^{*,}, 5*S*^{*)}-3-Acetyl-7-{4-[3-(4-*tert*-butyl-2-methylphenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid methylphenethylamide formate salt

10 Synthesized according to typical procedures F and E from bicyclononene U3 and 4-*tert*-butyl-2-methylphenol. LC-MS: R_t = 0.93, ES+: 608.25.

Example 107

15 (*rac.*)-(1*R*^{*,}, 5*S*^{*)}-3-Acetyl-7-{4-[3-(3,4-dichlorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid methylphenethylamide formate salt

20 Synthesized according to typical procedures F and E from bicyclononene U3 and 3,4-dichlorophenol. LC-MS: R_t = 0.84; ES+: 606.12.

Example 108

25 (*rac.*)-(1*R*^{*,}, 5*S*^{*)}-3-Acetyl-7-{4-[3-(4-bromo-3-methylphenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid methylphenethylamide formate salt

Synthesized according to typical procedures F and E from bicyclononene U3 and 4-bromo-3-methylphenol. LC-MS: R_t = 0.85; ES+: 630.11.

30

Example 109

(rac.)-(1R, 5S*)-3-Acetyl-7-{4-[3-(3,4-dimethylphenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid methylphenethylamide formate salt*

5 Synthesized according to typical procedures F and E from bicyclononene U3 and
3,4-dimethylphenol. LC-MS: R_t = 0.82; ES+: 566.20.

Example 110

10 *(rac.)-(1R*, 5S*)-3-Acetyl-7-{4-[3-(3,5-dichlorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid methylphenethylamide formate salt*

15 Synthesized according to typical procedures F and E from bicyclononene U3 and
3,5-dichlorophenol. LC-MS: R_t = 0.87; ES+: 606.13.

Example 111

20 *(rac.)-(1R*, 5S*)-3-Acetyl-7-{4-[3-(3,5-dimethylphenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid methylphenethylamide formate salt*

Synthesized according to typical procedures F and E from bicyclononene U3 and
3,5-dimethylphenol. LC-MS: R_t = 0.82; ES+: 566.21.

25

Example 112

30 *(rac.)-(1R*, 5S*)-3-Acetyl-7-{4-[3-(2-chloro-4,5-dimethylphenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid methylphenethylamide formate salt*

Synthesized according to typical procedures F and E from bicyclononene U3 and 2-chloro-4,5-dimethylphenol. LC-MS: R_t = 0.85; ES+: 600.18.

Example 113

5

(*rac.*)-(IR*, 5S*)-3-Acetyl-7-{4-[3-(2,3,5-trimethylphenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid methylphenethylamide formate salt

10 Synthesized according to typical procedures F and E from bicyclononene U3 and 2,3,5-trimethylphenol. LC-MS: R_t = 0.86; ES+: 580.23.

Example 114

15

(*rac.*)-Acetic acid (IR*, 5S*)-2-(7-{4-[3-(2-chlorophenoxy)propyl]phenyl}-6-[(2-(2-chlorophenyl)ethyl)methylcarbamoyl]-3,9-diazabicyclo[3.3.1]non-6-en-3-yl)-2-oxoethyl ester formate salt

20

Synthesized according to typical procedures A and E from bicyclononene AC and acetic acid chlorocarbonylmethyl ester. LC-MS: R_t = 1.01; ES+: 664.14.

Example 115

25

(*rac.*)-(IR*, 5S*)-7-{4-[3-(2-Chlorophenoxy)propyl]phenyl}-3-(2-cyanoacetyl)-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid [2-(2-chlorophenyl)ethyl]-methylamide formate salt

30

Example 116

Synthesized according to typical procedures K and E from bicyclononene AC and cyanoacetic acid. LC-MS: R_t = 1.02; ES+: 631.13.

(*rac.*)-(1*R**, 5*S**)-3-(2-Acetylaminoacetyl)-7-{4-[3-(2-chlorophenoxy)propyl]-phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid [2-(2-chlorophenyl)ethyl]methylamide formate salt

- 5 Synthesized according to typical procedures K and E from bicyclononene AC and acetylaminoacetic acid. LC-MS: R_t = 0.96; ES+: 663.14.

Example 117

- 10 1:1 Mixture of (1*R*, 5*S*)-3-((4*S*)-2-acetylamino-4-methylpentanoyl)-7-{4-[3-(2-chlorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid [2-(2-chlorophenyl)ethyl]methylamide formate salt and (1*S*, 5*R*)-3-((4*S*)-2-acetyl-amino-4-methylpentanoyl)-7-{4-[3-(2-chlorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid [2-(2-chlorophenyl)ethyl]methylamide formate salt
15

Synthesized according to typical procedures K and E from bicyclononene AC and acetyl leucine. LC-MS: R_t = 1.05; ES+: 719.19.

20 **Example 118**

(*rac.*)-(1*R**, 5*S**)-3-Acetyl-7-{4-[3-(2-chlorophenoxy)propyl]phenyl}-3,9-di-azabicyclo[3.3.1]non-6-ene-6-carboxylic acid [2-(2-chlorophenyl)ethyl]-methylamide formate salt

- 25 Synthesized according to typical procedures G and E from bicyclononene T2 and [2-(2-chlorophenyl)ethyl]methylamine (Jaques B.; Wallace R. G., *Tetrahedron*, 1977, 33, 581). LC-MS: R_t = 0.90; ES+: 606.08.

30 **Example 119**

(*rac.*)-(1*R*^{*}, 5*S*^{*})-3-Acetyl-7-{4-[3-(2-chlorophenoxy)propyl]phenyl}-3,9-di-azabicyclo[3.3.1]non-6-ene-6-carboxylic acid (2-hydroxybenzyl)methylamide formate salt

- 5 Synthesized according to typical procedures G and E from bicyclononene T2 and 2-methylaminomethylphenol (Ross S. D., *et al.*; *J. Org. Chem.*, **1966**, *31*, 133). LC-MS: R_t = 0.83; ES+: 574.10.

Example 120

10

(*rac.*)-(1*R*^{*}, 5*S*^{*})-3-Acetyl-7-{4-[3-(2-chlorophenoxy)propyl]phenyl}-3,9-di-azabicyclo[3.3.1]non-6-ene-6-carboxylic acid (2-chlorobenzyl)cyclopropylamide formate salt

15

- Synthesized according to typical procedures G and E from bicyclononene T2 and (2-chlorobenzyl)cyclopropylamine. LC-MS: R_t = 0.92; ES+: 617.94.

Example 121

20

- 1:1 Mixture of (*rac.*)-(1*R*^{*}, 5*S*^{*})-3-acetyl-7-{4-[3-(2-chlorophenoxy)propyl]-phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid [(3*R*^{*})3-(2-chlorophenyl)butyl]methylamide formate salt and (*rac.*)-(1*R*^{*}, 5*S*^{*})-3-acetyl-7-{4-[3-(2-chlorophenoxy)propyl]-phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid [(3*S*^{*})3-(2-chlorophenyl)butyl]methylamide formate salt

25

- Synthesized according to typical procedures G and E from bicyclononene T2 and (*rac.*)-methyl(3-phenylbutyl)amine (Meyers A. I., *et al.*; *J. Am. Chem. Soc.*, **1982**, *104*, 877). LC-MS: R_t = 0.91; ES+: 600.13.

30 **Example 122**

(*rac.*)-(1*R*^{*}, 5*S*^{*})-3-Acetyl-7-{4-[3-(2-chlorophenoxy)propyl]phenyl}-3,9-di-azabicyclo[3.3.1]non-6-ene-6-carboxylic acid methyl-(4-phenylbutyl)amide formate salt

- 5 Synthesized according to typical procedures G and E from bicyclononene T2 and methyl(4-phenylbutyl)amine (Neale R. S., *et al.*; *J. Org. Chem.*, **1965**, *30*, 3683). LC-MS: R_t = 0.91; ES+: 600.20.

Example 123

- 10 **(*rac.*)-(1*R*^{*}, 5*S*^{*})-3-Acetyl-7-{4-[3-(2-chlorophenoxy)propyl]phenyl}-3,9-di-azabicyclo[3.3.1]non-6-ene-6-carboxylic acid methyl-(3-phenoxypropyl)amide formate salt**

- 15 Synthesized according to typical procedures G and E from bicyclononene T2 and methyl(3-phenoxypropyl)amine. LC-MS: R_t = 0.88; ES+: 602.09.

Example 124

- 20 **(*rac.*)-(1*R*^{*}, 5*S*^{*})-3-Acetyl-7-{4-[3-(2-chlorophenoxy)propyl]phenyl}-3,9-di-azabicyclo[3.3.1]non-6-ene-6-carboxylic acid methyl-(4-phenylpentyl)amide formate salt**

- 25 Synthesized according to typical procedures G and E from bicyclononene T2 and methyl(5-phenylpentyl)amine (Neale R. S., *et al.*; *J. Org. Chem.*, **1965**, *30*, 3683). LC-MS: R_t = 0.95; ES+: 614.12.

Example 125

- 30 **(*rac.*)-(1*R*^{*}, 5*S*^{*})-3-Acetyl-7-{4-[3-(2-chlorophenoxy)propyl]phenyl}-3,9-di-azabicyclo[3.3.1]non-6-ene-6-carboxylic acid (3-benzo[1,3]dioxol-5-ylpropyl)-methylamide formate salt**

Synthesized according to typical procedures G and E from bicyclononene T2 and (3-benzo[1,3]dioxol-5-ylpropyl)methylamine (Dallacker, *et al.*; *Chem. Ber.*, 1971, 104, 2517). LC-MS: R_t = 0.86; ES+: 630.10.

5

Example 126

(*rac.*)-(IR*, 5S*)-3-Acetyl-7-{4-[3-(2-chlorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid [2-(4-methoxyphenoxy)ethyl]-methylamide formate salt

Synthesized according to typical procedures G and E from bicyclononene T2 and [2-(4-methoxyphenoxy)ethyl]methylamine. LC-MS: R_t = 0.84; ES+: 618.03.

15 **Example 127**

(*rac.*)-(IR*, 5S*)-3-Acetyl-7-{4-[3-(2-chlorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid [2-(4-chlorophenoxy)ethyl]-methylamide formate salt

20

Synthesized according to typical procedures G and E from bicyclononene T2 and [2-(4-chlorophenoxy)ethyl]methylamine. LC-MS: R_t = 0.90; ES+: 622.03.

Example 128

25

(*rac.*)-(IR*, 5S*)-3-Acetyl-7-{4-[3-(2-chlorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid methyl-(2-p-tolyloxyethyl)-amide formate salt

30 Synthesized according to typical procedures G and E from bicyclononene T2 and (2-p-tolyloxyethyl)methylamine. LC-MS: R_t = 0.89; ES+: 602.08.

Example 129

(rac.)-(1R, 5S*)-3-Acetyl-7-{4-[3-(2-chlorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid diethylamide formate salt*

5

Synthesized according to typical procedures G and E from bicyclononene T2 and diethylamine. LC-MS: R_t = 0.79; ES+: 510.06.

Example 130

10

(rac.)-(1R, 5S*)-3-Acetyl-7-{4-[3-(2-chlorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid methyl-(2-pyridin-2-ylethyl)-amide formate salt*

15

Synthesized according to typical procedures G and E from bicyclononene T2 and ethyl(2-pyridin-2-ylethyl)amine. LC-MS: R_t = 0.89; ES+: 602.08.

Example 131

20

(1S, 5R)-3-Acetyl-7-{4-[2-(2,3,6-trifluorophenoxy)ethyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid methylphenethylamide

Synthesized according to typical procedures F and E from bicyclononene BJ and 2,3,6-trifluorophenol. LC-MS: R_t = 0.94; ES+: 592.19. ee = 80%.

25

Example 132

1:1-Mixture of (rac.)-(1R, 5S*)-3-acetyl-7-{4-[3-(2-bromo-5-fluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid cyclopropyl-[2-((2R*)-2,3-dihydroxypropyl)benzyl]amide trifluoroacetate salt and (rac.)-(1R*, 5S*)-3-acetyl-7-{4-[3-(2-bromo-5-fluorophenoxy)-propyl]phenyl}-*

3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid cyclopropyl-[2-((2S*)-2,3-dihydroxypropyl)benzyl]amide trifluoroacetate salt

Synthesized according to typical procedure E from bicyclononene AT. LC-MS:
5 R_t = 3.99; ES+: 720.49.

Example 133

(*rac.*)-(1*R**, 5*S**)-3-Acetyl-7-{4-[3-(2-bromo-5-fluorophenoxy)propyl]phenyl}-
10 3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid cyclopropyl-[2-(2-hydroxyethyl)benzyl]amide trifluoroacetate salt

Synthesized according to typical procedure E from bicyclononene AV. LC-MS:
R_t = 3.94; ES+: 692.77.

15

Example 134

(*rac.*)-(1*R**, 5*S**)-3-Acetyl-7-{4-[3-(2-bromo-5-fluorophenoxy)propyl]phenyl}-
15 3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid benzylcyclopropylamide
20 formate salt

Synthesized according to typical procedures H and E from bicyclononene AJ4 and benzylcyclopropylamine. LC-MS: R_t = 0.89 ES+: 646.41.

25 **Example 135**

(*rac.*)-(1*R**, 5*S**)-3-Acetyl-7-{4-[3-(2-bromo-5-fluoro-phenoxy)propyl]-phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid (2-chlorobenzyl)-ethlamide formate salt

30

Synthesized according to typical procedures H and E from bicyclononene AJ4 and (2-chlorobenzyl)ethylamine. LC-MS: R_t = 0.91 ES+: 668.44.

Example 136

5 *(rac.)-(1R*, 5S*)-3-Acetyl-7-{4-[3-(2-bromo-5-fluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid cyclopropyl-(2-fluorobenzyl)amide formate salt*

Synthesized according to typical procedures H and E from bicyclononene AJ4 and (2-fluorobenzyl)cyclopropylamine. LC-MS: R_t = 0.90 ES+: 664.46.

10

Example 137

15 *(rac.)-(1R*, 5S*)-3-Acetyl-7-{4-[3-(2-bromo-5-fluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid cyclopropyl-(2-methylbenzyl)amide formate salt*

Synthesized according to typical procedures H and E from bicyclononene AJ4 and (2-methylbenzyl)cyclopropylamine. LC-MS: R_t = 0.92 ES+: 660.47.

20 **Example 138**

(rac.)-(1R, 5S*)-3-Acetyl-7-{4-[3-(2-bromo-5-fluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid cyclopropyl-[2-(4-methoxyphenoxy)ethyl]amide formate salt*

25

Synthesized according to typical procedures H and E from bicyclononene AJ4 and cyclopropyl[2-(4-methoxyphenoxy)ethyl]amine. LC-MS: R_t = 0.90 ES+: 706.44.

30 **Example 139**

(rac.)-(IR, 5S*)-3-Acetyl-7-{4-[3-(2-bromo-5-fluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid cyclopropyl-(2-m-tolyloxyethyl)amide formate salt*

- 5 Synthesized according to typical procedures H and E from bicyclononene AJ4 and cyclopropyl[2-(3-methylphenoxy)ethyl]amine. LC-MS: R_t = 0.93 ES+: 690.47.

Example 140

10

(rac.)-(IR, 5S*)-3-Acetyl-7-{4-[3-(2-bromo-5-fluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid cyclopropyl-[2-(3,4-dimethylphenoxy)ethyl]amide formate salt*

15

- Synthesized according to typical procedures H and E from bicyclononene AJ4 and cyclopropyl[2-(3,4-dimethylphenoxy)ethyl]amine. LC-MS: R_t = 0.94 ES+: 704.48.

Example 141

20

(rac.)-(IR, 5S*)-3-Acetyl-7-{4-[3-(2-bromo-5-fluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid cyclopropylphenethylamide formate salt*

25

- Synthesized according to typical procedures H and E from bicyclononene AJ4 and cyclopropylphenethylamine. LC-MS: R_t = 0.90 ES+: 660.50.

Example 142

30

- (rac.)-(IR*, 5S*)-3-Acetyl-7-{4-[3-(2-bromo-5-fluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid [2-(2-chlorophenyl)ethyl]-cyclopropylamide formate salt*

Synthesized according to typical procedures H and E from bicyclononene **AJ4** and [2-(2-chlorophenyl)ethyl]cyclopropylamine. LC-MS: R_t = 0.92 ES+: 694.44.

5 **Example 143**

(*rac.*)-(IR*, 5S*)-3-Acetyl-7-{4-[3-(2-bromo-5-fluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid cyclopropyl-[2-(2,3-difluorophenyl)ethyl]amide formate salt

10

Synthesized according to typical procedures H and E from bicyclononene **AJ4** and [2-(2,3-difluorophenyl)ethyl]cyclopropylamine. LC-MS: R_t = 0.91 ES+: 696.47.

15 **Example 144**

(*rac.*)-(IR*, 5S*)-3-Acetyl-7-{4-[3-(2-bromo-5-fluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid cyclopropyl-[2-(4-fluorophenyl)ethyl]amide formate salt

20

Synthesized according to typical procedures H and E from bicyclononene **AJ4** and [2-(4-fluorophenyl)ethyl]cyclopropylamine. LC-MS: R_t = 0.91 ES+: 678.53.

Example 145

25

(*rac.*)-(IR*, 5S*)-3-Acetyl-7-{4-[3-(2-bromo-5-fluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid cyclopropyl-(2-*o*-tolylethyl)amide formate salt

30

Synthesized according to typical procedures H and E from bicyclononene **AJ4** and [2-(2-methylphenyl)ethyl]cyclopropylamine. LC-MS: R_t = 0.92 ES+: 674.55.

Example 146

1:1-Mixture of (*rac.*)-(1*R*^{*}, 5*S*^{*})-3-Acetyl-7-{4-[3-(2-bromo-5-fluorophenoxy)-propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid ((2*R*^{*})-2-hydroxy-2-phenylethyl)methylamide formate salt and (*rac.*)-(1*R*^{*}, 5*S*^{*})-3-acetyl-7-{4-[3-(2-bromo-5-fluorophenoxy)-propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid ((2*S*^{*})-2-hydroxy-2-phenylethyl)methylamide formate salt

10 Synthesized according to typical procedures G and E from bicyclononene AJ4 and α-(methylaminomethyl)benzyl alcohol. LC-MS: R_t = 0.85 ES+: 650.49.

Example 147

15 (*rac.*)-(1*R*^{*}, 5*S*^{*})-3-Acetyl-7-{4-[3-(2-bromo-5-fluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid cyclopropyl-(3-trifluoromethylbenzyl)amide formate salt

20 Synthesized according to typical procedures H and E from bicyclononene AJ4 and (3-trifluoromethylbenzyl)cyclopropylamine. LC-MS: R_t = 0.93 ES+: 714.40.

Example 148

25 (*rac.*)-(1*R*^{*}, 5*S*^{*})-3-Acetyl-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid cyclopropyl-(2-*o*-tolylethyl)amide formate salt

Synthesized according to typical procedures H and E from bicyclononene AJ4 and [2-(2-methylphenyl)ethyl]cyclopropylamine. LC-MS: R_t = 0.90 ES+: 632.51.

Example 149

(*rac.*)-(IR*, 5S*)-7-{4-[3-(2,3,6-Trifluorophenoxy)propylphenyl]-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid (2-chlorobenzyl)cyclopropylamide

Synthesized according to typical procedure E from bicyclononene **AL2** and
5 purification by FC. LC-MS: R_t = 0.84 ES+: 596.30.

Example 150

(*rac.*)-5-((IR*, 5S*)-6-[(2-Chlorobenzyl)cyclopropylcarbamoyl]-7-{4-[3-(2,3,6-trifluorophenoxy)propylphenyl]-3,9-diazabicyclo[3.3.1]non-6-en-3-yl)-5-oxo-pentanoic acid formate salt

Synthesized according to typical procedures K and E from bicyclononene **AL2** and glutaric anhydride. LC-MS: R_t = 0.87 ES+: 710.42.
15

Example 151

(*rac.*)-(IR*, 5S*)-6-[(2-Chlorobenzyl)cyclopropylcarbamoyl]-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-9-carboxylic acid 2,2,2-trichloro-1,1-dimethylethyl ester hydrochloride salt

Synthesized according to typical procedure L from bicyclononene **AK2** and. LC-MS: R_t = 1.06 ES+: 798.34.

25 **Example 152**

(*rac.*)-(IR*, 5S*)-6-[(2-Chlorobenzyl)ethylcarbamoyl]-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-9-carboxylic acid 2,2,2-trichloro-1,1-dimethylethyl ester hydrochloride salt

30 Synthesized according to typical procedure L from bicyclononene **AK4** and. LC-MS: R_t = 1.19 ES+: 788.25.

Example 153

(*rac.*)-(1*R*^{*}, 5*S*^{*})-6-[(2-Fluorobenzyl)cyclopropylcarbamoyl]-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-9-carboxylic acid 2,2,2-trichloro-1,1-dimethylethyl ester hydrochloride salt

Synthesized according to typical procedure L from bicyclononene AK5. LC-MS:

R_t = 1.18 ES+: 784.27.

10

Example 154

(*rac.*)-(1*R*^{*}, 5*S*^{*})-6-[(2-Methylbenzyl)cyclopropylcarbamoyl]-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-9-carboxylic acid 2,2,2-trichloro-1,1-dimethylethyl ester hydrochloride salt

Synthesized according to typical procedure L from bicyclononene AK7. LC-MS:

R_t = 1.21 ES+: 780.29.

20

Example 155

(*rac.*)-(1*R*^{*}, 5*S*^{*})-6-[Cyclopropyl-(2-*o*-tolylethyl)carbamoyl]-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-9-carboxylic acid 2,2,2-trichloro-1,1-dimethylethyl ester hydrochloride salt

25

Synthesized according to typical procedure L from bicyclononene AK16. LC-MS: R_t = 1.21 ES+: 794.30.

30

Example 156

(rac.)-(1R, 5S*)-6-[Cyclopropyl-(3,5-dimethoxybenzyl)carbamoyl]-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-9-carboxylic acid 2,2,2-trichloro-1,1-dimethylethyl ester hydrochloride salt*

5 Synthesized according to typical procedure L from bicyclononene AK17. LC-MS: R_t = 1.16.

Example 157

10 *(rac.)-(1R*, 5S*)-6-[Cyclopropyl-(2-p-tolyethyl)carbamoyl]-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-9-carboxylic acid 2,2,2-trichloro-1,1-dimethylethyl ester hydrochloride salt*

Synthesized according to typical procedure L from bicyclononene AK18. LC-MS: R_t = 1.22 ES+: 794.30.

Example 158

20 *(rac.)-5-((1R*, 5S*)-6-[(2-Chlorobenzyl)cyclopropylcarbamoyl]-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-en-3-yl)-2,2-dimethyl-5-oxo-pentanoic acid formate salt*

Synthesized according to typical procedures K and E from bicyclononene AL2 and 2,2-dimethylglutaric anhydride. LC-MS: R_t = 0.91 ES+: 738.52.

25

Example 159

30 *(rac.)-5-((1R*, 5S*)-6-[(2-Chlorobenzyl)cyclopropylcarbamoyl]-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-en-3-yl)-5-oxo-pentanoic acid methyl ester formate salt*

Synthesized according to typical procedures A and E from bicyclononene AL2 and glutaric acid monomethylester chloride. LC-MS: R_t = 0.93 ES+: 724.49.

Example 160

5

1:1-Mixture of (*1R*, *5S*)-3-((*1S*, *4R*)-4-hydroxypyrrolidine-2-carbonyl)-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid (2-chlorobenzyl)cyclopropylamide formate salt and (*1S*, *5R*)-3-((*1S*, *4R*)-4-hydroxypyrrolidine-2-carbonyl)-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid (2-chlorobenzyl)cyclopropylamide formate salt

10

Synthesized according to typical procedures G, E and L from bicyclononene AL2 and BOC-L-hydroxyproline. LC-MS: R_t = 0.80 ES+: 709.38.

15

Example 161

20

(*rac.*)-(1*R*^{*}, 5*S*^{*})-3-(4-Carbamoylbutyryl)-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid (2-chlorobenzyl)cyclopropylamide formate salt

25

Synthesized according to typical procedures G and E from bicyclononene AL2 and 4-carbamoylbutyric acid. LC-MS: R_t = 0.86 ES+: 709.38.

30

Example 162

(*rac.*)-(1*R*^{*}, 5*S*^{*})-3-(4-Carbamoylbutyryl)-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid benzylcyclopropylamide formate salt

35

Synthesized according to typical procedures G and E from bicyclononene AL3 and 4-carbamoylbutyric acid. LC-MS: R_t = 0.84 ES+: 675.49.

Example 163

(*rac.*)-(1*R*^{*}, 5*S*^{*})-3-(4-Carbamoylbutyryl)-7-{4-[3-(2,3,6-trifluorophenoxy)-propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid (2-chlorobenzyl)ethylamide formate salt

Synthesized according to typical procedures G and E from bicyclononene AL4 and 4-carbamoylbutyric acid. LC-MS: R_t = 0.86 ES+: 697.40.

10

Example 164

(*rac.*)-(1*R*^{*}, 5*S*^{*})-3-(4-Carbamoylbutyryl)-7-{4-[3-(2,3,6-trifluorophenoxy)-propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid cyclopropyl-(2-fluorobenzyl)amide formate salt

Synthesized according to typical procedures G and E from bicyclononene AL5 and 4-carbamoylbutyric acid. LC-MS: R_t = 0.85 ES+: 693.43.

20 **Example 165**

(*rac.*)-(1*R*^{*}, 5*S*^{*})-3-(4-Carbamoylbutyryl)-7-{4-[3-(2,3,6-trifluorophenoxy)-propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid cyclopropyl-(3-trifluoromethylbenzyl)amide formate salt

25

Synthesized according to typical procedures G and E from bicyclononene AL6 and 4-carbamoylbutyric acid. LC-MS: R_t = 0.88 ES+: 743.41.

Example 166

30

(*rac.*)-(1*R*^{*}, 5*S*^{*})-3-(4-Carbamoylbutyryl)-7-{4-[3-(2,3,6-trifluorophenoxy)-propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid cyclo-propyl-(2-methylbenzyl)amide formate salt

- 5 Synthesized according to typical procedures G and E from bicyclononene AL7 and 4-carbamoylbutyric acid. LC-MS: R_t = 0.86 ES+: 689.47.

Example 167

- 10 (*rac.*)-(1*R*^{*}, 5*S*^{*})-3-(4-Carbamoylbutyryl)-7-{4-[3-(2,3,6-trifluorophenoxy)-propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid cyclo-propyl-[2-(4-methoxyphenoxy)ethyl]amide formate salt

15 Synthesized according to typical procedures G and E from bicyclononene AL8 and 4-carbamoylbutyric acid. LC-MS: R_t = 0.85 ES+: 735.47.

Example 168

- 20 (*rac.*)-(1*R*^{*}, 5*S*^{*})-3-(4-Carbamoylbutyryl)-7-{4-[3-(2,3,6-trifluorophenoxy)-propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid cyclo-propyl-[2-(3-methoxyphenoxy)ethyl]amide formate salt

25 Synthesized according to typical procedures G and E from bicyclononene AL9 and 4-carbamoylbutyric acid. LC-MS: R_t = 0.86 ES+: 735.47.

Example 169

- 30 (*rac.*)-(1*R*^{*}, 5*S*^{*})-3-(4-Carbamoylbutyryl)-7-{4-[3-(2,3,6-trifluorophenoxy)-propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid cyclo-propyl-(2-*m*-tolyloxyethyl)amide formate salt

Synthesized according to typical procedures G and E from bicyclononene AL10 and 4-carbamoylbutyric acid. LC-MS: R_t = 0.87 ES+: 719.46.

Example 170

5

(*rac.*)-(IR*, 5S*)-3-(4-Carbamoylbutyryl)-7-{4-[3-(2,3,6-trifluorophenoxy)-propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid cyclo-propyl-[2-(3,4-dimethylphenoxy)ethyl]amide formate salt

10 Synthesized according to typical procedures G and E from bicyclononene AL11 and 4-carbamoylbutyric acid. LC-MS: R_t = 0.89 ES+: 733.49.

Example 171

15 (*rac.*)-(IR*, 5S*)-3-(4-Carbamoylbutyryl)-7-{4-[3-(2,3,6-trifluorophenoxy)-propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid cyclo-propylphenethylamide formate salt

20 Synthesized according to typical procedures G and E from bicyclononene AL12 and 4-carbamoylbutyric acid. LC-MS: R_t = 0.85 ES+: 689.48.

Example 172

25 (*rac.*)-(IR*, 5S*)-3-(4-Carbamoylbutyryl)-7-{4-[3-(2,3,6-trifluorophenoxy)-propyl]phenyl}-3,9-diaza-bicyclo[3.3.1]non-6-ene-6-carboxylic acid [2-(2-chloro-phenyl)ethyl]cyclopropylamide formate salt

Synthesized according to typical procedures G and E from bicyclononene AL13 and 4-carbamoylbutyric acid. LC-MS: R_t = 0.87 ES+: 723.43.

30

Example 173

(rac.)-(1R, 5S*)-3-(4-Carbamoylbutyryl)-7-{4-[3-(2,3,6-trifluorophenoxy)-propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid cyclo-propyl-[2-(2,3-difluorophenyl)ethyl]amide formate salt*

- 5 Synthesized according to typical procedures G and E from bicyclononene **AL14** and 4-carbamoylbutyric acid. LC-MS: R_t = 0.86 ES+: 725.45.

Example 174

- 10 *(rac.)-(1R*, 5S*)-3-(4-Carbamoylbutyryl)-7-{4-[3-(2,3,6-trifluorophenoxy)-propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid cyclo-propyl-[2-(4-fluorophenyl)ethyl]amide*

15 Synthesized according to typical procedures G and E from bicyclononene **AL15** and 4-carbamoylbutyric acid. LC-MS: R_t = 0.86 ES+: 707.44.

Example 175

- 20 *(rac.)-(1R*, 5S*)-3-(4-Carbamoylbutyryl)-7-{4-[3-(2,3,6-trifluorophenoxy)-propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid cyclopropyl-(2-*o*-tolylethyl)amide formate salt*

25 Synthesized according to typical procedures G and E from bicyclononene **AL16** and 4-carbamoylbutyric acid. LC-MS: R_t = 0.87 ES+: 703.47.

25

Example 176

- 30 *(rac.)-(1R*, 5S*)-3-(4-Carbamoylbutyryl)-7-{4-[3-(2,3,6-trifluorophenoxy)-propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid cyclo-propyl-(3,5-dimethoxybenzyl)amide formate salt*

Synthesized according to typical procedures G and E from bicyclononene **AL17** and 4-carbamoylbutyric acid. LC-MS: R_t = 0.85 ES+: 735.47.

Example 177

5

(*rac.*)-(1*R**, 5*S**)-3-(4-Carbamoylbutyryl)-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid cyclo-propyl-(2-*p*-tolylethyl)amide formate salt

10 Synthesized according to typical procedures G and E from bicyclononene **AL18** and 4-carbamoylbutyric acid. LC-MS: R_t = 0.87 ES+: 703.46.

Example 178

15 1:1-Mixture of (*IR*, 5*S*)-3-((1*S*, 4*R*)-4-hydroxypyrrolidine-2-carbonyl)-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid benzylcyclopropylamide formate salt and (*IS*, 5*R*)-3-((1*S*, 4*R*)-4-hydroxypyrrolidine-2-carbonyl)-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid benzyl-cyclopropylamide formate salt
20

Synthesized according to typical procedures G, E and L from bicyclononene **AL3** and BOC-L-hydroxyproline. LC-MS: R_t = 0.78 ES+: 675.47.

25 **Example 179**

1:1-Mixture of (*IR*, 5*S*)-3-((1*S*, 4*R*)-4-hydroxypyrrolidine-2-carbonyl)-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid (2-chlorobenzyl)ethylamide formate salt and (*IS*, 5*R*)-3-((1*S*, 4*R*)-4-hydroxypyrrolidine-2-carbonyl)-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid (2-chlorobenzyl)ethylamide formate salt
30

Synthesized according to typical procedures G, E and L from bicyclononene **AL4** and BOC-L-hydroxyproline. LC-MS: R_t = 0.79 ES+: 697.40.

5 **Example 180**

1:1-Mixture of (*1R, 5S*)-3-((*1S, 4R*)-4-hydroxypyrrolidine-2-carbonyl)-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid cyclopropyl-(2-fluorobenzyl)amide formate salt and (*1S, 5R*)-3-((*1S, 4R*)-4-hydroxypyrrolidine-2-carbonyl)-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid cyclopropyl-(2-fluorobenzyl)amide formate salt

10 Synthesized according to typical procedures G, E and L from bicyclononene **AL5** and BOC-L-hydroxyproline. LC-MS: R_t = 0.78 ES+: 693.44.

Example 181

1:1-Mixture of (*1R, 5S*)-3-((*1S, 4R*)-4-hydroxypyrrolidine-2-carbonyl)-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid cyclopropyl-(3-trifluoromethylbenzyl)amide formate salt and (*1S, 5R*)-3-((*1S, 4R*)-4-hydroxypyrrolidine-2-carbonyl)-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid cyclopropyl-(3-trifluoromethylbenzyl)amide formate salt

15 Synthesized according to typical procedures G, E and L from bicyclononene **AL6** and BOC-L-hydroxyproline. LC-MS: R_t = 0.81 ES+: 743.42.

Example 182

30 1:1-Mixture of (*1R, 5S*)-3-((*1S, 4R*)-4-hydroxypyrrolidine-2-carbonyl)-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-

6-carboxylic acid cyclopropyl-(2-methylbenzyl)amide formate salt and (*1S*,
5R)-3-((*1S*, *4R*)-4-hydroxypyrrolidine-2-carbonyl)-7-{4-[3-(2,3,6-trifluoro-
phenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid
cyclopropyl-(2-methylenzyl)amide formate salt

5

Synthesized according to typical procedures G, E and L from bicyclononene AL7
and BOC-L-hydroxyproline. LC-MS: R_t = 0.79 ES+: 689.47.

Example 183

10

1:1-Mixture of (*1R*, *5S*)-3-((*1S*, *4R*)-4-hydroxypyrrolidine-2-carbonyl)-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-
6-carboxylic acid cyclopropyl-(2-*m*-tolyloxyethyl)amide formate salt and (*1S*,
5R)-3-((*1R* *4S*)-4-hydroxypyrrolidine-2-carbonyl)-7-{4-[3-(2,3,6-trifluoro-
phenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid
cyclopropyl-(2-*m*-tolyloxyethyl)amide formate salt

15

Synthesized according to typical procedures G, E and L from bicyclononene
AL10 and BOC-L-hydroxyproline. LC-MS: R_t = 0.81 ES+: 719.45.

20

Example 184

1:1-Mixture of (*1R*, *5S*)-3-((*1S*, *4R*)-4-hydroxypyrrolidine-2-carbonyl)-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-
6-carboxylic acid cyclopropyl-(3,5-dimethoxybenzyl)amide fromate salt and
25 (*1S*, *5R*)-3-((*1S*, *4R*)-4-hydroxypyrrolidine-2-carbonyl)-7-{4-[3-(2,3,6-
trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carbo-
xylic acid cyclopropyl-(3,5-dimethoxybenzyl)amide formate salt

25

30 Synthesized according to typical procedures G, E and L from bicyclononene
AL17 and BOC-L-hydroxyproline. LC-MS: R_t = 0.78 ES+: 735.48.

Example 185

1:1-Mixture of (*1R, 5S*)-3-((*1S, 4R*)-hydroxypyrrolidine-2-carbonyl)-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid cyclopropyl-(2-*p*-tolylethyl)amide formate salt and (*1S, 5R*)-3-((*1S, 4R*)-hydroxypyrrolidine-2-carbonyl)-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid cyclopropyl-(2-*p*-tolylethyl)-amide formate salt

10 Synthesized according to typical procedures G, E and L from bicyclononene AL18 and BOC-L-hydroxypyroline. LC-MS: R_t = 0.81 ES+: 703.48.

Example 186

15 1:1-Mixture of (*rac.*)-(3*R*^{*})-5-((*1R*^{*}, *5S*^{*})-6-[(2-chlorobenzyl)cyclopropylcarbamoyl]-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-en-3-yl)-3-hydroxy-5-oxopentanoic acid formate salt and (*rac.*)-(3*S*^{*})-5-((*1R*^{*}, *5S*^{*})-6-[(2-chlorobenzyl)cyclopropylcarbamoyl]-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-en-3-yl)-3-hydroxy-5-oxopentanoic acid formate salt

20 Synthesized according to typical procedures K, E and L from bicyclononene AL2 and 3-(*tert*-butyldimethylsilyloxy)glutaric anhydride. LC-MS: R_t = 0.86 ES+: 726.40.

25

Example 187

1:1-Mixture of (*rac.*)-(3*R*^{*})-5-((*1R*^{*}, *5S*^{*})-6-(benzylcyclopropylcarbamoyl)-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-en-3-yl)-3-hydroxy-5-oxopentanoic acid formate salt and (*rac.*)-(3*S*^{*})-5-((*1R*^{*}, *5S*^{*})-6-(benzylcyclopropylcarbamoyl)-7-{4-[3-(2,3,6-trifluorophenoxy)

propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-en-3-yl)-3-hydroxy-5-oxo-pentanoic acid formate salt

Synthesized according to typical procedures K, E and L from bicyclononene AL3
5 and 3-(*tert*-butyldimethylsilyloxy)glutaric anhydride. LC-MS: R_t = 0.84 ES+:
692.45

Example 188

10 **1:1-Mixture of (*rac.*)-(3*R*^{*})-5-((1*R*^{*}, 5*S*^{*})-6-[*(2-chlorobenzyl)ethylcarbamoyl*]-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo-[3.3.1]non-6-en-3-yl)-3-hydroxy-5-oxopentanoic acid formate salt and (*rac.*)-(3*S*^{*})-5-((1*R*^{*}, 5*S*^{*})-6-[*(2-chlorobenzyl)ethylcarbamoyl*]-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-en-3-yl)-3-**
15 **hydroxy-5-oxopentanoic acid formate salt**

Synthesized according to typical procedures K, E and L from bicyclononene AL4
and 3-(*tert*-butyldimethylsilyloxy)glutaric anhydride. LC-MS: R_t = 0.86 ES+:
714.38

20

Example 189

1:1-Mixture of (*rac.*)-(3*R*^{*})-5-((1*R*^{*}, 5*S*^{*})-6-[cyclopropyl-(2-fluorobenzyl)-carbamoyl]-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo-[3.3.1]non-6-en-3-yl)-3-hydroxy-5-oxopentanoic acid formate salt and (*rac.*)-(3*S*^{*})-5-((1*R*^{*}, 5*S*^{*})-6-[cyclopropyl-(2-fluorobenzyl)carbamoyl]-7-{4-[3-(2,3,6-trifluoro-phenoxy)propyl]phenyl}-3,9-diazabicyclo-[3.3.1]non-6-en-3-yl)-3-hydroxy-5-oxopentanoic acid formate salt

30 **Synthesized according to typical procedures K, E and L from bicyclononene AL5 and 3-(*tert*-butyldimethylsilyloxy)glutaric anhydride. LC-MS: R_t = 0.84 ES+: 710.42**

Example 190

1:1-Mixture of (*rac.*)-(3*R*^{*})-5-((1*R*^{*}, 5*S*^{*})-6-[cyclopropyl-(3-trifluoromethylbenzyl)carbamoyl]-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-en-3-yl)-3-hydroxy-5-oxopentanoic acid formate salt and (*rac.*)-(3*S*^{*})-5-((1*R*^{*}, 5*S*^{*})-6-[cyclopropyl-(3-trifluoromethylbenzyl)carbamoyl]-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-en-3-yl)-3-hydroxy-5-oxopentanoic acid formate salt

10

Synthesized according to typical procedures K, E and L from bicyclononene AL6 and 3-(*tert*-butyldimethylsilyloxy)glutaric anhydride. LC-MS: R_t = 0.88 ES+: 760.39

15 **Example 191**

1:1-Mixture of (*rac.*)-(3*R*^{*})-5-((1*R*^{*}, 5*S*^{*})-6-[cyclopropyl-(2-methylbenzyl)carbamoyl]-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-en-3-yl)-3-hydroxy-5-oxopentanoic acid formate salt and (*rac.*)-(3*S*^{*})-5-((1*R*^{*}, 5*S*^{*})-6-[cyclopropyl-(2-methylbenzyl)carbamoyl]-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-en-3-yl)-3-hydroxy-5-oxopentanoic acid formate salt

Synthesized according to typical procedures K, E and L from bicyclononene AL7 and 3-(*tert*-butyldimethylsilyloxy)glutaric anhydride. LC-MS: R_t = 0.86 ES+: 706.44

Example 192

30 1:1-Mixture of (*rac.*)-(3*R*^{*})-5-((1*R*^{*}, 5*S*^{*})-6-[cyclopropyl-[2-(4-methoxyphenoxy)ethyl]carbamoyl]-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-en-3-yl)-3-hydroxy-5-oxopentanoic acid formate

salt and (*rac.*)-(3*S**)-5-((*IR**, 5*S**)-6-{cyclopropyl-[2-(4-methoxyphenoxy)-ethyl]carbamoyl}-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-en-3-yl)-3-hydroxy-5-oxopentanoic acid formate salt

- 5 Synthesized according to typical procedures K, E and L from bicyclononene AL8 and 3-(*tert*-butyldimethylsilyloxy)glutaric anhydride. LC-MS: R_t = 0.85 ES+:
752.43

Example 193

- 10 1:1-Mixture of (*rac.*)-(3*R**)-5-((*IR**, 5*S**)-6-{cyclopropyl-[2-(3-methoxyphenoxy)ethyl]carbamoyl}-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-en-3-yl)-3-hydroxy-5-oxopentanoic acid formate salt and (*rac.*)-(3*S**)-5-((*IR**, 5*S**)-6-{cyclopropyl-[2-(3-methoxyphenoxy)-ethyl]carbamoyl}-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-en-3-yl)-3-hydroxy-5-oxopentanoic acid formate salt

- 15 Synthesized according to typical procedures K, E and L from bicyclononene AL9 and 3-(*tert*-butyldimethylsilyloxy)glutaric anhydride. LC-MS: R_t = 0.85 ES+:
20 752.44

Example 194

- 25 1:1-Mixture of (*rac.*)-(3*R**)-5-((*IR**, 5*S**)-6-[cyclopropyl-(2-*m*-tolyloxyethyl)carbamoyl]-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo-[3.3.1]non-6-en-3-yl)-3-hydroxy-5-oxopentanoic acid formate salt and (*rac.*)-(3*S**)-5-((*IR**, 5*S**)-6-[cyclopropyl-(2-*m*-tolyloxyethyl)carbamoyl]-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo-[3.3.1]non-6-en-3-yl)-3-hydroxy-5-oxopentanoic acid formate salt .

Synthesized according to typical procedures K, E and L from bicyclononene AL10 and 3-(*tert*-butyldimethylsilyloxy)glutaric anhydride. LC-MS: R_t = 0.87
ES+: 736.45.

5 **Example 195**

10 **1:1-Mixture of (*rac.*)-(3*R*^{*})-5-((1*R*^{*}, 5*S*^{*})-6-{cyclopropyl-[2-(3,4-dimethylphenoxy)ethyl]carbamoyl}-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-en-3-yl)-3-hydroxy-5-oxopentanoic acid formate salt and (*rac.*)-(3*S*^{*})-5-((1*R*^{*}, 5*S*^{*})-6-{cyclopropyl-[2-(3,4-dimethylphenoxy)ethyl]carbamoyl}-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-en-3-yl)-3-hydroxy-5-oxopentanoic acid formate salt**

15 Synthesized according to typical procedures K, E and L from bicyclononene AL11 and 3-(*tert*-butyldimethylsilyloxy)glutaric anhydride. LC-MS: R_t = 0.89
ES+: 750.47

Example 196

20 **1:1-Mixture of (*rac.*)-(3*R*^{*})-5-((1*R*^{*}, 5*S*^{*})-(6-(cyclopropylphenethylcarbamoyl)-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-en-3-yl)-3-hydroxy-5-oxopentanoic acid formate salt and (*rac.*)-(3*S*^{*})-5-((1*R*^{*}, 5*S*^{*})-5-(6-(cyclopropylphenethylcarbamoyl)-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-en-3-yl)-3-**

25 **hydroxy-5-oxopentanoic acid formate salt**

Synthesized according to typical procedures K, E and L from bicyclononene AL12 and 3-(*tert*-butyldimethylsilyloxy)glutaric anhydride. LC-MS: R_t = 0.85
ES+: 706.43

1:1-Mixture of (*rac.*)-(3*R*^{*})-5-((*IR*^{*}, 5*S*^{*})-6-{[2-(2-chlorophenyl)ethyl]-cyclopropylcarbamoyl}-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-en-3-yl)-3-hydroxy-5-oxopentanoic acid formate salt and (*rac.*)-(3*S*^{*})-5-((*IR*^{*}, 5*S*^{*})-6-{[2-(2-chlorophenyl)ethyl]cyclopropyl-carbamoyl}-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-en-3-yl)-3-hydroxy-5-oxopentanoic acid formate salt

Synthesized according to typical procedures K, E and L from bicyclononene AL13 and 3-(*tert*-butyldimethylsilyloxy)glutaric anhydride. LC-MS: R_t = 0.87
10 ES+: 740.40

Example 198

1:1-Mixture of (*rac.*)-(3*R*^{*})-5-((*IR*^{*}, 5*S*^{*})-6-{cyclopropyl-[2-(2,3-difluoro-phenyl)ethyl]carbamoyl}-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-en-3-yl)-3-hydroxy-5-oxopentanoic acid formate salt and (*rac.*)-(3*S*^{*})-5-((*IR*^{*}, 5*S*^{*})-6-{cyclopropyl-[2-(2,3-difluorophenyl)ethyl]-carbamoyl}-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-en-3-yl)-3-hydroxy-5-oxopentanoic acid formate salt

Synthesized according to typical procedures K, E and L from bicyclononene AL14 and 3-(*tert*-butyldimethylsilyloxy)glutaric anhydride. LC-MS: R_t = 0.86
20 ES+: 742.42

Example 199

1:1-Mixture of (*rac.*)-(3*R*^{*})-5-((*IR*^{*}, 5*S*^{*})-6-{cyclopropyl-[2-(4-fluorophenyl)-ethyl]carbamoyl}-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-en-3-yl)-3-hydroxy-5-oxopentanoic acid formate salt and (*rac.*)-(3*S*^{*})-5-((*IR*^{*}, 5*S*^{*})-6-{cyclopropyl-[2-(4-fluorophenyl)ethyl]-carbamoyl}-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-en-3-yl)-3-hydroxy-5-oxopentanoic acid formate salt

Synthesized according to typical procedures K, E and L from bicyclononene **AL15** and 3-(*tert*-butyldimethylsilyloxy)glutaric anhydride. LC-MS: R_t = 0.86
ES+: 724.43.

5

Example 200

1:1-Mixture of (*rac.*)-(3*R*^{*})-5-((1*R*^{*}, 5*S*^{*})-6-[cyclopropyl-(2-*o*-tolylethyl)carbamoyl]-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-en-3-yl)-3-hydroxy-5-oxopentanoic acid formate salt and (*rac.*)-(3*S*^{*})-5-((1*R*^{*}, 5*S*^{*})-6-[cyclopropyl-(2-*o*-tolylethyl)carbamoyl]-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-en-3-yl)-3-hydroxy-5-oxopentanoic acid formate salt

15 Synthesized according to typical procedures K, E and L from bicyclononene **AL16** and 3-(*tert*-butyldimethylsilyloxy)glutaric anhydride. LC-MS: R_t = 0.87
ES+: 720.45

20

1:1-Mixture of (*rac.*)-(3*R*^{*})-5-((1*R*^{*}, 5*S*^{*})-6-[cyclopropyl-(3,5-dimethoxybenzyl)carbamoyl]-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-en-3-yl)-3-hydroxy-5-oxopentanoic acid formate salt and (*rac.*)-(3*S*^{*})-5-((1*R*^{*}, 5*S*^{*})-6-[cyclopropyl-(3,5-dimethoxybenzyl)carbamoyl]-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-en-3-yl)-3-hydroxy-5-oxopentanoic acid formate salt

30 ES+: 752.42

Example 202

1:1-Mixture of (*rac.*)-(3*R**)-5-((1*R**, 5*S**)-6-[cyclopropyl-(2-*p*-tolylethyl)-carbamoyl]-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo-[3.3.1]non-6-en-3-yl)-3-hydroxy-5-oxopentanoic acid formate salt and (*rac.*)-(3*R**)-5-((1*R**, 5*S**)-6-[cyclopropyl-(2-*p*-tolylethyl)carbamoyl]-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo-[3.3.1]non-6-en-3-yl)-3-hydroxy-5-oxopentanoic acid formate salt

Synthesized according to typical procedures K, E and L from bicyclononene
10 **AL18** and 3-(*tert*-butyldimethylsilyloxy)glutaric anhydride. LC-MS: R_t = 0.87
ES+: 720.43

Example 203

15 (*rac.*)-(1*R**, 5*S**)-3-Acetyl-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid benzylcyclopropylamide formate salt

Synthesized according to typical procedures A and E from bicyclononene **AL3**
20 and acetyl chloride. LC-MS: R_t = 0.88 ES+: 604.53

Example 204

25 (*rac.*)-(1*R**, 5*S**)-3-Acetyl-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid (2-chlorobenzyl)ethylamide formate salt

Synthesized according to typical procedures A and E from bicyclononene **AL4**
and acetyl chloride. LC-MS: R_t = 0.90 ES+: 626.48

Example 205

(*rac.*)-(1*R*^{*, 5*S*})-3-Acetyl-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid cyclopropyl(2-fluorobenzyl)-amide formate salt

5 Synthesized according to typical procedures A and E from bicyclononene AL5 and acetyl chloride. LC-MS: R_t = 0.89 ES+: 622.53

Example 206

10 **(*rac.*)-(1*R*^{*, 5*S*})-3-Acetyl-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid cyclopropyl-(2-methylbenzyl)-amide formate salt**

15 Synthesized according to typical procedures A and E from bicyclononene AL7 and acetyl chloride. LC-MS: R_t = 0.90 ES+: 618.54

Example 207

20 **(*rac.*)-(1*R*^{*, 5*S*})-3-Acetyl-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid cyclopropyl[2-(4-methoxy-phenoxy)ethyl]amide formate salt**

25 Synthesized according to typical procedures A and E from bicyclononene AL8 and acetyl chloride. LC-MS: R_t = 0.89 ES+: 664.54

Example 208

30 **(*rac.*)-(1*R*^{*, 5*S*})-3-Acetyl-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid cyclopropyl[2-(3-methoxy-phenoxy)ethyl]amide formate salt**

Synthesized according to typical procedures A and E from bicyclononene **AL9** and acetyl chloride. LC-MS: $R_t = 0.89$ ES+: 664.53

Example 209

5

(*rac.*)-(1*R**, 5*S**)-3-Acetyl-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid cyclopropyl[2-(3-methylphenoxy)ethyl]amide formate salt

10 Synthesized according to typical procedures A and E from bicyclononene **AL10** and acetyl chloride. LC-MS: $R_t = 0.91$ ES+: 648.53

Example 210

15 (*rac.*)-(1*R**, 5*S**)-3-Acetyl-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid cyclopropylphenethylamide formate salt

20 Synthesized according to typical procedures A and E from bicyclononene **AL12** and acetyl chloride. LC-MS: $R_t = 0.89$ ES+: 618.54

Example 211

25 (*rac.*)-(1*R**, 5*S**)-3-Acetyl-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid cyclopropyl(3,5-dimethoxybenzyl)amide formate salt

Synthesized according to typical procedures A and E from bicyclononene **AL17** and acetyl chloride. LC-MS: $R_t = 0.88$ ES+: 664.55

30

Example 212

(*rac.*)-(1*R*^{*, 5*S*})-3-Acetyl-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid cyclopropyl(2-*p*-tolylethyl)amide formate salt

5 Synthesized according to typical procedures A and E from bicyclononene AL18 and acetyl chloride. LC-MS: R_t = 0.91 ES+: 632.54

Example 213

10 (*rac.*)-5-((1*R*^{*, 5*S*})-6-(Benzylcyclopropylcarbamoyl)-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-en-3-yl)-5-oxopentanoic acid formate salt

15 Synthesized according to typical procedures K and E from bicyclononene AL3 and glutaric anhydride. LC-MS: R_t = 0.86 ES+: 676.54

Example 214

20 (*rac.*)-5-((1*R*^{*, 5*S*})-6-[(2-Chlorobenzyl)ethylcarbamoyl]-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-en-3-yl)-5-oxopentanoic acid formate salt

25 Synthesized according to typical procedures K and E from bicyclononene AL4 and glutaric anhydride. LC-MS: R_t = 0.88 ES+: 698.46

Example 215

30 (*rac.*)-5-((1*R*^{*, 5*S*})-6-[Cyclopropyl-(2-fluorobenzyl)carbamoyl]-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-en-3-yl)-5-oxopentanoic acid formate salt

Synthesized according to typical procedures K and E from bicyclononene **AL5** and glutaric anhydride. LC-MS: R_t = 0.86 ES+: 694.51

Example 216

5

(*rac.*)-5-((1*R*^{*, 5*S*^{*})-6-[Cyclopropyl(3-trifluoromethylbenzyl)carbamoyl]-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-en-3-yl)-5-oxopentanoic acid formate salt}

10 Synthesized according to typical procedures K and E from bicyclononene **AL6** and glutaric anhydride. LC-MS: R_t = 0.89 ES+: 744.51

Example 217

15

(*rac.*)-5-((1*R*^{*, 5*S*^{*})-6-[Cyclopropyl(2-methylbenzyl)carbamoyl]-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-en-3-yl)-5-oxo-pentanoic acid formate salt}

20

Synthesized according to typical procedures K and E from bicyclononene **AL7** and glutaric anhydride. LC-MS: R_t = 0.88 ES+: 690.54

Example 218

25

(*rac.*)-5-((1*R*^{*, 5*S*^{*})-6-{Cyclopropyl[2-(4-methoxyphenoxy)ethyl]carbamoyl}-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-en-3-yl)-5-oxopentanoic acid formate salt}

Synthesized according to typical procedures K and E from bicyclononene **AL8** and glutaric anhydride. LC-MS: R_t = 0.87 ES+: 736.54

30

Example 219

(rac.)-5-((1R, 5S*)-6-{Cyclopropyl[2-(3-methoxyphenoxy)ethyl]carbamoyl}-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-en-3-yl)-5-oxopentanoic acid formate salt*

- 5 Synthesized according to typical procedures K and E from bicyclononene **AL9** and glutaric anhydride. LC-MS: R_t = 0.87 ES+: 736.55

Example 220

- 10 *(rac.)-5-((1R*, 5S*)-6-{Cyclopropyl[2-(3-methylphenoxy)ethyl]carbamoyl}-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-en-3-yl)-5-oxopentanoic acid formate salt*

15 Synthesized according to typical procedures K and E from bicyclononene **AL10** and glutaric anhydride. LC-MS: R_t = 0.89 ES+: 720.53

Example 221

- 20 *(rac.)-5-((1R*, 5S*)-6-{Cyclopropyl-[2-(3,4-dimethylphenoxy)ethyl]-carbamoyl}-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo-[3.3.1]non-6-en-3-yl)-5-oxopentanoic acid formate salt*

25 Synthesized according to typical procedures K and E from bicyclononene **AL11** and glutaric anhydride. LC-MS: R_t = 0.90 ES+: 734.57

Example 222

- 30 *(rac.)-5-((1R*, 5S*)-6-(Cyclopropylphenethylcarbamoyl)-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-en-3-yl)-5-oxopentanoic acid formate salt*

Synthesized according to typical procedures K and E from bicyclononene **AL12** and glutaric anhydride. LC-MS: R_t = 0.87 ES+: 690.52

Example 223

5

(*rac.*)-5-((*1R*^{*,} *5S*^{*})-6-{[2-(2-Chlorophenyl)ethyl]cyclopropylcarbamoyl}-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-en-3-yl)-5-oxopentanoic acid formate salt

10 Synthesized according to typical procedures K and E from bicyclononene **AL13** and glutaric anhydride. LC-MS: R_t = 0.89 ES+: 724.49

Example 224

15 (*rac.*)-5-((*1R*^{*,} *5S*^{*})-6-{Cyclopropyl-[2-(2,3-difluorophenyl)ethyl]carbamoyl}-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-en-3-yl)-5-oxopentanoic acid formate salt

20 Synthesized according to typical procedures K and E from bicyclononene **AL14** and glutaric anhydride. LC-MS: R_t = 0.88ES+: 726.51

Example 225

25 (*rac.*)-5-((*1R*^{*,} *5S*^{*})-6-{Cyclopropyl[2-(4-fluorophenyl)ethyl]carbamoyl}-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-en-3-yl)-5-oxopentanoic acid formate salt

Synthesized according to typical procedures K and E from bicyclononene **AL15** and glutaric anhydride. LC-MS: R_t = 0.87 ES+: 708.50

30

Example 226

(*rac.*)-5-((*1R*^{*}, *5S*^{*})-6-{Cyclopropyl[2-(2-methylphenyl)ethyl]carbamoyl}-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-en-3-yl)-5-oxopentanoic acid formate salt

5 Synthesized according to typical procedures K and E from bicyclononene **AL16** and glutaric anhydride. LC-MS: R_t = 0.88 ES+: 704.54

Example 227

10 **(*rac.*)-5-((*1R*^{*}, *5S*^{*})-6-[Cyclopropyl(3,5-dimethoxybenzyl)carbamoyl]-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-en-3-yl)-5-oxopentanoic acid formate salt**

15 Synthesized according to typical procedures K and E from bicyclononene **AL17** and glutaric anhydride. LC-MS: R_t = 0.86 ES+: 736.55

Example 228

20 **(*rac.*)-5-((*1R*^{*}, *5S*^{*})-6-[Cyclopropyl(2-p-tolylethyl)carbamoyl]-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-en-3-yl)-5-oxopentanoic acid formate salt**

25 Synthesized according to typical procedures K and E from bicyclononene **AL18** and glutaric anhydride. LC-MS: R_t = 0.88 ES+: 704.54

Example 229

30 **(*rac.*)-5-((*1R*^{*}, *5S*^{*})-6-(Benzylcyclopropylcarbamoyl)-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-en-3-yl)-5-oxopen-**

tanoic acid methyl ester formate salt

Synthesized according to typical procedures A and E from bicyclononene **AL3** and glutaric acid monomethylester chloride. LC-MS: R_t = 0.91 ES+: 690.55

Example 230

5

(*rac.*)-5-((*1R*^{*}, *5S*^{*})-6-[*(2-Chlorobenzyl)ethylcarbamoyl*]-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-en-3-yl)-5-oxopentanoic acid methyl ester formate salt

10 Synthesized according to typical procedures A and E from bicyclononene **AL4** and glutaric acid monomethylester chloride. LC-MS: R_t = 0.92 ES+: 712.49

Example 231

15 (*rac.*)-5-((*1R*^{*}, *5S*^{*})-6-[*Cyclopropyl-(2-fluorobenzyl)carbamoyl*]-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-en-3-yl)-5-oxopentanoic acid methyl ester formate salt

20 Synthesized according to typical procedures A and E from bicyclononene **AL5** and glutaric acid monomethylester chloride. LC-MS: R_t = 0.91 ES+: 708.51

Example 232

25 (*rac.*)-5-((*1R*^{*}, *5S*^{*})-6-[*Cyclopropyl(3-trifluoromethylbenzyl)carbamoyl*]-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-en-3-yl)-5-oxopentanoic acid methyl ester formate salt

Synthesized according to typical procedures A and E from bicyclononene **AL6** and glutaric acid monomethylester chloride. LC-MS: R_t = 0.94 ES+: 758.51

30

Example 233

(*rac.*)-5-((*1R*^{*}, *5S*^{*})-6-[Cyclopropyl(2-methylbenzyl)carbamoyl]-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-en-3-yl)-5-oxopentanoic acid methyl ester formate salt

5 Synthesized according to typical procedures A and E from bicyclononene AL7 and glutaric acid monomethylester chloride. LC-MS: R_t = 0.92 ES+: 704.54

Example 234

10 **(*rac.*)-5-((*1R*^{*}, *5S*^{*})-6-[Cyclopropyl[2-(4-methoxyphenoxy)ethyl]carbamoyl]-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-en-3-yl)-5-oxopentanoic acid methyl ester formate salt**

15 Synthesized according to typical procedures A and E from bicyclononene AL8 and glutaric acid monomethylester chloride. LC-MS: R_t = 0.92 ES+: 750.54

Example 235

20 **(*rac.*)-5-((*1R*^{*}, *5S*^{*})-6-[Cyclopropyl[2-(3-methoxyphenoxy)ethyl]carbamoyl]-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-en-3-yl)-5-oxopentanoic acid methyl ester formate salt**

Synthesized according to typical procedures A and E from bicyclononene AL9 and glutaric acid monomethylester chloride. LC-MS: R_t = 0.92 ES+: 750.56

25

Example 236

30 **(*rac.*)-5-((*1R*^{*}, *5S*^{*})-6-[Cyclopropyl[2-(3-methylphenoxy)ethyl]carbamoyl]-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-en-3-yl)-5-oxopentanoic acid methyl ester formate salt**

Synthesized according to typical procedures A and E from bicyclononene **AL10** and glutaric acid monomethylester chloride. LC-MS: R_t = 0.93 ES+: 734.58

Example 237

5

(*rac.*)-5-((*1R*^{*}, *5S*^{*})-6-{Cyclopropyl[2-(3,4-dimethylphenoxy)ethyl]-carbamoyl}-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo-[3.3.1]non-6-en-3-yl)-5-oxopentanoic acid methyl ester formate salt

10 Synthesized according to typical procedures A and E from bicyclononene **AL11** and glutaric acid monomethylester chloride. LC-MS: R_t = 0.95 ES+: 748.57

Example 238

15 (*rac.*)-5-((*1R*^{*}, *5S*^{*})-6-(Cyclopropylphenethylcarbamoyl)-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-en-3-yl)-5-oxopentanoic acid methyl ester formate salt

20 Synthesized according to typical procedures A and E from bicyclononene **AL12** and glutaric acid monomethylester chloride. LC-MS: R_t = 0.91 ES+: 704.55

Example 239

25 (*rac.*)-5-((*1R*^{*}, *5S*^{*})-6-{[2-(2-Chlorophenyl)ethyl]cyclopropylcarbamoyl}-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-en-3-yl)-5-oxopentanoic acid methyl ester formate salt

30 Synthesized according to typical procedures A and E from bicyclononene **AL13** and glutaric acid monomethylester chloride. LC-MS: R_t = 0.93 ES+: 738.53

Example 240

(*rac.*)-5-((*1R*^{*}, *5S*^{*})-6-{Cyclopropyl[2-(2,3-difluorophenyl)ethyl]carbamoyl}-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-en-3-yl)-5-oxopentanoic acid methyl ester formate salt

5 Synthesized according to typical procedures A and E from bicyclononene **AL14** and glutaric acid monomethylester chloride. LC-MS: R_t = 0.92 ES+: 740.54

Example 241

10 **(*rac.*)-5-((*1R*^{*}, *5S*^{*})-6-{Cyclopropyl[2-(4-fluorophenyl)ethyl]carbamoyl}-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-en-3-yl)-5-oxopentanoic acid methyl ester formate salt**

15 Synthesized according to typical procedures A and E from bicyclononene **AL15** and glutaric acid monomethylester chloride. LC-MS: R_t = 0.92 ES+: 722.54

Example 242

20 **(*rac.*)-5-((*1R*^{*}, *5S*^{*})-6-{Cyclopropyl[2-(2-methylphenyl)ethyl]carbamoyl}-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-en-3-yl)-5-oxopentanoic acid methyl ester formate salt**

25 Synthesized according to typical procedures A and E from bicyclononene **AL16** and glutaric acid monomethylester chloride. LC-MS: R_t = 0.93 ES+: 718.56

Example 243

30 **(*rac.*)-5-((*1R*^{*}, *5S*^{*})-6-[Cyclopropyl-(3,5-dimethoxybenzyl)carbamoyl]-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-en-3-yl)-5-oxopentanoic acid methyl ester formate salt**

Synthesized according to typical procedures A and E from bicyclononene AL17 and glutaric acid monomethylester chloride. LC-MS: R_t = 0.91 ES+: 750.55

Example 244

5

(*rac.*)-5-((*1R*^{*}, *5S*^{*})-6-{Cyclopropyl[2-(4-methylphenyl)ethyl]carbamoyl}-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-en-3-yl)-5-oxopentanoic acid methyl ester formate salt

10 Synthesized according to typical procedures A and E from bicyclononene AL18 and glutaric acid monomethylester chloride. LC-MS: R_t = 0.93 ES+: 718.56

Example 245

15 (*rac.*)-5-((*1R*^{*}, *5S*^{*})-6-(Benzylcyclopropylcarbamoyl)-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-en-3-yl)-2,2-dimethyl-5-oxopentanoic acid methyl ester formate salt

20 Synthesized according to typical procedures K and E from bicyclononene AL3 and 2,2-dimethylglutaric anhydride. LC-MS: R_t = 0.90 ES+: 704.53

Example 246

25 (*rac.*)-5-((*1R*^{*}, *5S*^{*})-6-[(2-Chlorobenzyl)ethylcarbamoyl]-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-en-3-yl)-2,2-dimethyl-5-oxopentanoic acid formate salt

Synthesized according to typical procedures K and E from bicyclononene AL4 and 2,2-dimethylglutaric anhydride. LC-MS: R_t = 0.91 ES+: 726.53

30

Example 247

(*rac.*)-5-((*1R, *5S**)-6-[Cyclopropyl-(2-fluorobenzyl)carbamoyl]-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-en-3-yl)-2,2-dimethyl-5-oxo-pentanoic formate salt**

- 5 Synthesized according to typical procedures K and E from bicyclononene AL5 and 2,2-dimethylglutaric anhydride. LC-MS: R_t = 0.90 ES+: 722.54

Example 248

10 **(*rac.*)-5-((*1R**, *5S**)-6-[Cyclopropyl-(3-trifluoromethylbenzyl)carbamoyl]-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-en-3-yl)-2,2-dimethyl-5-oxopentanoic acid formate salt**

- 15 Synthesized according to typical procedures K and E from bicyclononene AL6 and 2,2-dimethylglutaric anhydride. LC-MS: R_t = 0.93 ES+: 772.51

Example 249

20 **(*rac.*)-5-((*1R**, *5S**)-6-[Cyclopropyl-(2-methylbenzyl)carbamoyl]-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-en-3-yl)-2,2-dimethyl-5-oxopentanoic acid formate salt**

- 25 Synthesized according to typical procedures K and E from bicyclononene AL7 and 2,2-dimethylglutaric anhydride. LC-MS: R_t = 0.91 ES+: 718.57

25

Example 250

30 **(*rac.*)-5-((*1R**, *5S**)-6-{Cyclopropyl-[2-(4-methoxyphenoxy)ethyl]carbamoyl}-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-en-3-yl)-2,2-dimethyl-5-oxopentanoic acid formate salt**

Synthesized according to typical procedures K and E from bicyclononene AL8 and 2,2-dimethylglutaric anhydride. LC-MS: R_t = 0.90 ES+: 764.55

Example 251

5

(*rac.*)-5-((1*R*^{*}, 5*S*^{*})-6-{Cyclopropyl-[2-(3-methoxyphenoxy)ethyl]carbamoyl}-7-{4-[3-(2,3,6-trifluoro-phenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-en-3-yl)-2,2-dimethyl-5-oxopentanoic acid formate salt

10 Synthesized according to typical procedures K and E from bicyclononene AL9 and 2,2-dimethylglutaric anhydride. LC-MS: R_t = 0.90 ES+: 764.54

Example 252

15 (*rac.*)-5-((1*R*^{*}, 5*S*^{*})-6-{Cyclopropyl-[2-(3-methylphenoxy)ethyl]carbamoyl}-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-en-3-yl)-2,2-dimethyl-5-oxopentanoic acid formate salt

20 Synthesized according to typical procedures K and E from bicyclononene AL10 and 2,2-dimethylglutaric anhydride. LC-MS: R_t = 0.92 ES+: 748.58

Example 253

25 (*rac.*)-5-((1*R*^{*}, 5*S*^{*})-6-{Cyclopropyl-[2-(3,4-dimethylphenoxy)ethyl]-carbamoyl}-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo-[3.3.1]non-6-en-3-yl)-2,2-dimethyl-5-oxopentanoic acid formate salt

Synthesized according to typical procedures K and E from bicyclononene AL11 and 2,2-dimethylglutaric anhydride. LC-MS: R_t = 0.93 ES+: 762.58

30

Example 254

(*rac.*)-5-((*1R*^{*}, *5S*^{*})-6-(Cyclopropylphenethylcarbamoyl)-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-en-3-yl)-2,2-dimethyl-5-oxo-pentanoic acid formate salt

- 5 Synthesized according to typical procedures K and E from bicyclononene AL12 and 2,2-dimethylglutaric anhydride. LC-MS: R_t = 0.90 ES+: 718.56

Example 255

- 10 (*rac.*)-5-((*1R*^{*}, *5S*^{*})-6-{[2-(2-Chlorophenyl)ethyl]cyclopropylcarbamoyl}-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-en-3-yl)-2,2-dimethyl-5-oxopentanoic acid formate salt

Synthesized according to typical procedures K and E from bicyclononene AL13
15 and 2,2-dimethylglutaric anhydride. LC-MS: R_t = 0.92 ES+: 752.50

Example 256

- 20 (*rac.*)-5-((*1R*^{*}, *5S*^{*})-6-{[2-(2,3-Difluorophenyl)ethyl]cyclopropylcarbamoyl}-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-en-3-yl)-2,2-dimethyl-5-oxopentanoic acid formate salt

Synthesized according to typical procedures K and E from bicyclononene AL14
25 and 2,2-dimethylglutaric anhydride. LC-MS: R_t = 0.91 ES+: 754.53

Example 257

- 30 (*rac.*)-5-((*1R*^{*}, *5S*^{*})-6-{[2-(4-Fluorophenyl)ethyl]cyclopropylcarbamoyl}-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-en-3-yl)-2,2-dimethyl-5-oxopentanoic acid formate salt

Synthesized according to typical procedures K and E from bicyclononene **AL15** and 2,2-dimethylglutaric anhydride. LC-MS: R_t = 0.91 ES+: 736.56

Example 258

5

(*rac.*)-5-((*1R*^{*}, *5S*^{*})-6-{[2-(2-Methylphenyl)ethyl]cyclopropylcarbamoyl}-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-en-3-yl)-2,2-dimethyl-5-oxopentanoic acid formate salt

10 Synthesized according to typical procedures K and E from bicyclononene **AL16** and 2,2-dimethylglutaric anhydride. LC-MS: R_t = 0.92 ES+: 732.59

Example 259

15 (*rac.*)-5-((*1R*^{*}, *5S*^{*})-6-[Cyclopropyl-(3,5-dimethoxybenzyl)carbamoyl]-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-en-3-yl)-2,2-dimethyl-5-oxopentanoic acid formate salt

20 Synthesized according to typical procedures K and E from bicyclononene **AL17** and 2,2-dimethylglutaric anhydride. LC-MS: R_t = 0.90 ES+: 764.54

Example 260

25 (*rac.*)-5-((*1R*^{*}, *5S*^{*})-6-{[2-(4-Methylphenyl)ethyl]cyclopropylcarbamoyl}-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-en-3-yl)-2,2-dimethyl-5-oxopentanoic acid formate salt

Synthesized according to typical procedures K and E from bicyclononene **AL18** and 2,2-dimethylglutaric anhydride. LC-MS: R_t = 0.92 ES+: 732.58

30

Example 261

1:1-Mixture of (*rac.*)-(2*R*^{*}, 3*S*^{*})-4-((1*R*^{*}, 5*S*^{*})-6-[(2-chlorobenzyl)ethylcarbamoyl]-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo-[3.3.1]non-6-en-3-yl)-2,3-dihydroxy-4-oxobutyric acid formate salt and (*rac.*)-(2*S*^{*}, 3*R*^{*})-4-((1*R*^{*}, 5*S*^{*})-6-[(2-chlorobenzyl)ethylcarbamoyl]-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo-[3.3.1]non-6-en-3-yl)-2,3-dihydroxy-4-oxobutyric acid formate salt

Synthesized according to typical procedures K and E from bicyclononene AL4 and *meso*-2,3-dihydroxysuccinic anhydride. LC-MS: R_t = 0.85 ES+: 716.45

10

Example 262

1:1-Mixture of (*rac.*)-(2*R*^{*}, 3*S*^{*})-4-((1*R*^{*}, 5*S*^{*})-6-[cyclopropyl-(2-fluorobenzyl)carbamoyl]-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-en-3-yl)-2,3-dihydroxy-4-oxobutyric acid formate salt and (*rac.*)-(2*S*^{*}, 3*R*^{*})-4-((1*R*^{*}, 5*S*^{*})-6-[cyclopropyl-(2-fluorobenzyl)carbamoyl]-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo-[3.3.1]non-6-en-3-yl)-2,3-dihydroxy-4-oxobutyric acid formate salt

20 Synthesized according to typical procedures K and E from bicyclononene AL5 and *meso*-2,3-dihydroxysuccinic anhydride. LC-MS: R_t = 0.83 ES+: 712.44

Example 263

25 1:1-Mixture of (*rac.*)-(2*R*^{*}, 3*S*^{*})-4-((1*R*^{*}, 5*S*^{*})-6-[Cyclopropyl-(3-trifluoromethylbenzyl)carbamoyl]-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo-[3.3.1]non-6-en-3-yl)-2,3-dihydroxy-4-oxobutyric acid formate salt and (*rac.*)-(2*S*^{*}, 3*R*^{*})-4-((1*R*^{*}, 5*S*^{*})-6-[cyclopropyl-(3-trifluoromethylbenzyl)carbamoyl]-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo-[3.3.1]non-6-en-3-yl)-2,3-dihydroxy-4-oxobutyric acid formate salt

Synthesized according to typical procedures K and E from bicyclononene AL6 and *meso*-2,3-dihydroxysuccinic anhydride. LC-MS: R_t = 0.87 ES+: 762.42

Example 264

5

1:1-Mixture of (*rac.*)-(2*R*^{*,} 3*S*^{*)}-4-((1*R*^{*,} 5*S*^{*)}-6-[cyclopropyl-(2-methylbenzyl)carbamoyl]-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-en-3-yl)-2,3-dihydroxy-4-oxobutyric acid formate salt and (*rac.*)-(2*S*^{*,} 3*R*^{*)}-4-((1*R*^{*,} 5*S*^{*)}-6-[cyclopropyl-(2-methylbenzyl)carbamoyl]-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo-[3.3.1]non-6-en-3-yl)-2,3-dihydroxy-4-oxobutyric acid formate salt

10

Synthesized according to typical procedures K and E from bicyclononene AL7 and *meso*-2,3-dihydroxysuccinic anhydride. LC-MS: R_t = 0.85 ES+: 708.46

15

Example 265

1:1-Mixture of (*rac.*)-(2*R*^{*,} 3*S*^{*)}-4-((1*R*^{*,} 5*S*^{*)}-6-[cyclopropyl-[2-(2,3-difluorophenyl)ethyl]carbamoyl]-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-en-3-yl)-2,3-dihydroxy-4-oxobutyric acid formate salt and (*rac.*)-(2*S*^{*,} 3*R*^{*)}-4-((1*R*^{*,} 5*S*^{*)}-6-[cyclopropyl-[2-(2,3-difluorophenyl)ethyl]carbamoyl]-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-en-3-yl)-2,3-dihydroxy-4-oxobutyric acid formate salt

20

Synthesized according to typical procedures K and E from bicyclononene AL14 and *meso*-2,3-dihydroxysuccinic anhydride. LC-MS: R_t = 0.85 ES+: 744.46

25

Example 266

30

1:1-Mixture of (*rac.*)-(2*R*^{*,} 3*S*^{*)}-4-((1*R*^{*,} 5*S*^{*)}-6-[cyclopropyl-[2-(2-methylphenyl)ethyl]carbamoyl]-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-

diazabicyclo[3.3.1]non-6-en-3-yl)-2,3-dihydroxy-4-oxobutyric acid formate salt and (*rac.*)-(2*S*^{*,} 3*R*^{*})-4-((1*R*^{*,} 5*S*^{*})-6-{cyclopropyl-[2-(2-methylphenyl)-ethyl]carbamoyl}-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-en-3-yl)-2,3-dihydroxy-4-oxobutyric acid formate salt

5

Synthesized according to typical procedures K and E from bicyclononene **AL16** and *meso*-2,3-dihydroxysuccinic anhydride. LC-MS: R_t = 0.86 ES+: 722.52

Example 267

10

1:1-Mixture of (*rac.*)-(2*R*^{*,} 3*S*^{*})-4-((1*R*^{*,} 5*S*^{*})-6-[cyclopropyl-(3,5-dimethoxybenzyl)carbamoyl]-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-en-3-yl)-2,3-dihydroxy-4-oxobutyric acid formate salt and (*rac.*)-(2*S*^{*,} 3*R*^{*})-4-((1*R*^{*,} 5*S*^{*})-6-[cyclopropyl-(3,5-dimethoxybenzyl)carbamoyl]-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-en-3-yl)-2,3-dihydroxy-4-oxobutyric acid formate salt

15

Synthesized according to typical procedures K and E from bicyclononene **AL17** and *meso*-2,3-dihydroxysuccinic anhydride. LC-MS: R_t = 0.84 ES+: 754.50

20

Example 268

1:1-Mixture of (*rac.*)-(2*R*^{*,} 3*S*^{*})-4-((1*R*^{*,} 5*S*^{*})-6-{cyclopropyl-[2-(4-methylphenyl)ethyl]carbamoyl}-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-en-3-yl)-2,3-dihydroxy-4-oxobutyric acid formate salt and (*rac.*)-(2*S*^{*,} 3*R*^{*})-4-((1*R*^{*,} 5*S*^{*})-6-{cyclopropyl-[2-(4-methylphenyl)ethyl]carbamoyl}-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-en-3-yl)-2,3-dihydroxy-4-oxobutyric acid formate salt

25

30 Synthesized according to typical procedures K and E from bicyclononene **AL18** and *meso*-2,3-dihydroxysuccinic anhydride. LC-MS: R_t = 0.86 ES+: 722.49

Example 269

- 1:1-Mixture of (*rac.*)-(3*R*^{*})-5-((*IR*^{*}, 5*S*^{*})-6-[(2-chlorobenzyl)cyclopropylcarbamoyl]-7-{4-[3-(2,3,5-trimethylphenoxy)ethyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-en-3-yl)-3-hydroxy-5-oxopentanoic acid formate salt and (*rac.*)-(3*S*^{*})-5-((*IR*^{*}, 5*S*^{*})-6-[(2-chlorobenzyl)cyclopropylcarbamoyl]-7-{4-[3-(2,3,5-trimethylphenoxy)ethyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-en-3-yl)-3-hydroxy-5-oxopentanoic acid formate salt
- 10 Synthesized according to typical procedures K, E and L from bicyclononene AL38 and 3-(*tert*-butyldimethylsilyloxy)glutaric anhydride. LC-MS: R_t = 0.88
ES+: 700.52

Example 270

- 15 1:1-Mixture of (*rac.*)-(3*R*^{*})-5-((*IR*^{*}, 5*S*^{*})-6-[(2-chlorobenzyl)ethylcarbamoyl]-7-{4-[3-(2,3,5-trimethylphenoxy)ethyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-en-3-yl)-3-hydroxy-5-oxopentanoic acid formate salt and (*rac.*)-(3*S*^{*})-5-((*IR*^{*}, 5*S*^{*})-6-[(2-chlorobenzyl)ethylcarbamoyl]-7-{4-[3-(2,3,5-trimethylphenoxy)ethyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-en-3-yl)-3-hydroxy-5-oxopentanoic acid formate salt

20 Synthesized according to typical procedures K, E and L from bicyclononene AL40 and 3-(*tert*-butyldimethylsilyloxy)glutaric anhydride. LC-MS: R_t = 0.88
25 ES+: 688.54

Example 271

- (*rac.*)-5-((*IR*^{*}, 5*S*^{*})-7-{4-[3-(2-Bromo-5-fluorophenoxy)propyl]phenyl}-6-[cyclopropyl-[2-(2-hydroxyethyl)benzyl]carbamoyl}-3,9-diazabicyclo[3.3.1]non-6-en-3-yl)-5-oxo-pentanoic acid formate salt

Synthesized according to typical procedures K and E from bicyclononene AL37 and glutaric anhydride. LC-MS: R_t = 0.86 ES+: 762.42

Example 272

5

(*rac.*)-5-((*1R*^{*}, *5S*^{*})-7-{4-[3-(2-Bromo-5-fluorophenoxy)propyl]phenyl}-6-{cyclopropyl-[2-(2-hydroxyethyl)benzyl]carbamoyl}-3,9-diazabicyclo[3.3.1]-non-6-en-3-yl)-5-oxo-pentanoic acid methyl ester formate salt

10 Synthesized according to typical procedures A and E from bicyclononene AL37 and glutaric acid monomethyl ester chloride. LC-MS: R_t = 0.88 ES+: 776.43

Example 273

15

(*rac.*)-(*1R*^{*}, *5S*^{*})-7-{4-[3-(2-Bromo-5-fluorophenoxy)propyl]phenyl}-3-(4-carbamoylbutyryl)-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid cyclopropyl-[2-(2-hydroxyethyl)benzyl]amide formate salt

20

Synthesized according to typical procedures G and E from bicyclononene AL37 and 4-carbamoylbutyric acid. LC-MS: R_t = 0.82 ES+: 761.45

Example 274

25

(*rac.*)-(*1R*^{*}, *5S*^{*})-3-Acetyl-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid cyclopropyl-[2-(2-hydroxyethyl)benzyl]amide formate salt

30

Synthesized according to typical procedures A and E from bicyclononene AL19 and acetyl chloride. LC-MS: R_t = 0.84 ES+: 648.50

Example 275

(*rac.*)-5-((*1R*^{*}, *5S*^{*})-6-{Cyclopropyl-[2-(2-hydroxyethyl)benzyl]carbamoyl}-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-en-3-yl)-5-oxo-pentanoic acid formate salt

- 5 Synthesized according to typical procedures K and E from bicyclononene **AL19** and glutaric anhydride. LC-MS: R_t = 0.85 ES+: 740.42

Example 276

10 (*rac.*)-5-((*1R*^{*}, *5S*^{*})-6-{Cyclopropyl-[2-(2-hydroxyethyl)benzyl]carbamoyl}-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-en-3-yl)-5-oxopentanoic acid methyl ester formate salt

- 15 Synthesized according to typical procedures A and E from bicyclononene **AL19** and glutaric acid monomethylester chloride. LC-MS: R_t = 0.87 ES+: 734.52

Example 277

20 (*rac.*)-5-((*1R*^{*}, *5S*^{*})-6-{Cyclopropyl-[2-(2-hydroxyethyl)benzyl]carbamoyl}-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-en-3-yl)-2,2-dimethyl-5-oxopentanoic acid formate salt

- Synthesized according to typical procedures K and E from bicyclononene **AL19** and 2,2-dimethylglutaric anhydride. LC-MS: R_t = 0.86 ES+: 748.52

25

Example 278

30 (*rac.*)-(*1R*^{*}, *5S*^{*})-3-(4-Carbamoylbutyryl)-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid cyclopropyl-[2-(2-hydroxyethyl)benzyl]amide formate salt

Synthesized according to typical procedures G and E from bicyclononene **AL19** and 4-carbamoylbutyric acid. LC-MS: R_t = 0.81; ES+: 719.52.

Example 279

5

(*rac.*)-5-{(1*R*^{*}, 5*S*^{*})-7-{4-[3-(2-Bromo-5-fluorophenoxy)propyl]phenyl}-6-[(2-chlorobenzyl)ethylcarbamoyl]-3,9-diazabicyclo[3.3.1]non-6-en-3-yl}-5-oxo-pentanoic acid formate salt

10 Synthesized according to typical procedures K and E from bicyclononene **AL22** and glutaric anhydride. LC-MS: R_t = 0.89; ES+: 740.38.

Example 280

15 (*rac.*)-5-{(1*R*^{*}, 5*S*^{*})-7-{4-[3-(2-Bromo-5-fluorophenoxy)propyl]phenyl}-6-[cyclopropyl-(2-fluorobenzyl)carbamoyl]-3,9-diazabicyclo[3.3.1]non-6-en-3-yl}-5-oxopentanoic acid formate salt

20 Synthesized according to typical procedures K and E from bicyclononene **AL23** and glutaric anhydride. LC-MS: R_t = 0.88; ES+: 736.41.

Example 281

25 (*rac.*)-5-{(1*R*^{*}, 5*S*^{*})-7-{4-[3-(2-Bromo-5-fluorophenoxy)propyl]phenyl}-6-[cyclopropyl-(3-trifluoromethylbenzyl)carbamoyl]-3,9-diazabicyclo[3.3.1]-non-6-en-3-yl}-5-oxopentanoic acid formate salt

Synthesized according to typical procedures K and E from bicyclononene **AL24** and glutaric anhydride. LC-MS: R_t = 0.91; ES+: 736.38.

30

Example 282

(*rac.*)-5-{(1*R*^{*}, 5*S*^{*})-7-[4-[3-(2-Bromo-5-fluorophenoxy)propyl]phenyl]-6-[cyclopropyl-(2-methylbenzyl)carbamoyl]-3,9-diazabicyclo[3.3.1]non-6-en-3-yl}-5-oxopentanoic acid formate salt

- 5 Synthesized according to typical procedures K and E from bicyclononene AL25 and glutaric anhydride. LC-MS: R_t = 0.89; ES+: 732.45.

Example 283

10 (*rac.*)-5-{(1*R*^{*}, 5*S*^{*})-7-[4-[3-(2-Bromo-5-fluorophenoxy)propyl]phenyl]-6-[cyclopropyl-[2-(4-methoxyphenoxy)ethyl]carbamoyl]-3,9-diazabicyclo-[3.3.1]non-6-en-3-yl}-5-oxopentanoic acid formate salt

- 15 Synthesized according to typical procedures K and E from bicyclononene AL26 and glutaric anhydride. LC-MS: R_t = 0.89; ES+: 778.41.

Example 284

20 (*rac.*)-5-{(1*R*^{*}, 5*S*^{*})-7-[4-[3-(2-Bromo-5-fluorophenoxy)propyl]phenyl]-6-[cyclopropyl-[2-(3-methylphenoxy)ethyl]carbamoyl]-3,9-diazabicyclo[3.3.1]-non-6-en-3-yl}-5-oxopentanoic acid formate salt

- 25 Synthesized according to typical procedures K and E from bicyclononene AL27 and glutaric anhydride. LC-MS: R_t = 0.91; ES+: 762.42.

Example 285

30 (*rac.*)-5-{(1*R*^{*}, 5*S*^{*})-7-[4-[3-(2-Bromo-5-fluorophenoxy)propyl]phenyl]-6-[cyclopropyl-[2-(3,4-dimethylphenoxy)ethyl]carbamoyl]-3,9-diazabicyclo-[3.3.1]non-6-en-3-yl}-5-oxopentanoic acid formate salt

Synthesized according to typical procedures K and E from bicyclononene **AL28** and glutaric anhydride. LC-MS: R_t = 0.92; ES+: 776.45.

Example 286

5

(*rac.*)-5-[(*1R*^{*}, *5S*^{*})-7-{4-[3-(2-Bromo-5-fluorophenoxy)propyl]phenyl}-6-(cyclopropylphenethylcarbamoyl)-3,9-diazabicyclo[3.3.1]non-6-en-3-yl]-5-oxopentanoic acid formate salt

10 Synthesized according to typical procedures K and E from bicyclononene **AL29** and glutaric anhydride. LC-MS: R_t = 0.89; ES+: 732.44.

Example 287

15 (*rac.*)-5-((*1R*^{*}, *5S*^{*})-7-{4-[3-(2-Bromo-5-fluorophenoxy)propyl]phenyl}-6-{[2-(2-chlorophenyl)ethyl]cyclopropylcarbamoyl}-3,9-diazabicyclo[3.3.1]non-6-en-3-yl)-5-oxopentanoic acid formate salt

Synthesized according to typical procedures K and E from bicyclononene **AL30**
20 and glutaric anhydride. LC-MS: R_t = 0.90; ES+: 766.35.

Example 288

25 (*rac.*)-5-((*1R*^{*}, *5S*^{*})-7-{4-[3-(2-Bromo-5-fluorophenoxy)propyl]phenyl}-6-{[2-(2,3-difluorophenyl)ethyl]cyclopropylcarbamoyl}-3,9-diazabicyclo[3.3.1]non-6-en-3-yl)-5-oxopentanoic acid formate salt

Synthesized according to typical procedures K and E from bicyclononene **AL31** and glutaric anhydride. LC-MS: R_t = 0.90; ES+: 768.35.

30

Example 289

(rac.)-5-((1R, 5S*)-7-{4-[3-(2-Bromo-5-fluorophenoxy)propyl]phenyl}-6-{[2-(4-fluorophenyl)ethyl]cyclopropylcarbamoyl}-3,9-diazabicyclo[3.3.1]non-6-en-3-yl)-5-oxopentanoic acid formate salt*

5 Synthesized according to typical procedures K and E from bicyclononene AL32 and glutaric anhydride. LC-MS: R_t = 0.89; ES+: 750.40.

Example 290

10 *(rac.)-5-((1R*, 5S*)-7-{4-[3-(2-Bromo-5-fluorophenoxy)propyl]phenyl}-6-{[2-(2-methylphenyl)ethyl]cyclopropylcarbamoyl}-3,9-diazabicyclo[3.3.1]non-6-en-3-yl)-5-oxopentanoic acid formate salt*

15 Synthesized according to typical procedures K and E from bicyclononene AL33 and glutaric anhydride. LC-MS: R_t = 0.90; ES+: 746.43.

Example 291

20 *(rac.)-5-((1R*, 5S*)-7-{4-[3-(2-Bromo-5-fluorophenoxy)propyl]phenyl}-6-[cyclopropyl-(3,5-dimethoxybenzyl)carbamoyl]-3,9-diazabicyclo[3.3.1]non-6-en-3-yl)-5-oxopentanoic acid formate salt*

25 Synthesized according to typical procedures K and E from bicyclononene AL35 and glutaric anhydride. LC-MS: R_t = 0.88; ES+: 778.40.

Example 292

30 *(rac.)-5-((1R*, 5S*)-7-{4-[3-(2-Bromo-5-fluorophenoxy)propyl]phenyl}-6-{[2-(4-methylphenyl)ethyl]cyclopropylcarbamoyl}-3,9-diazabicyclo[3.3.1]non-6-en-3-yl)-5-oxopentanoic acid formate salt*

Synthesized according to typical procedures K and E from bicyclononene AL36 and glutaric anhydride. LC-MS: R_t = 0.90; ES+: 746.43.

Example 293

5

(*rac.*)-5-{(1*R*^{*}, 5*S*^{*})-7-[4-[3-(2-Bromo-5-fluorophenoxy)propyl]phenyl]-6-[(2-chlorobenzyl)ethylcarbamoyl]-3,9-diazabicyclo[3.3.1]non-6-en-3-yl}-5-oxopentanoic acid methyl ester formate salt

10 Synthesized according to typical procedures A and E from bicyclononene AL22 and glutaric acid monomethylester chloride. LC-MS: R_t = 0.94; ES+: 754.37.

Example 294

15 (*rac.*)-5-{(1*R*^{*}, 5*S*^{*})-7-[4-[3-(2-Bromo-5-fluorophenoxy)propyl]phenyl]-6-[cyclopropyl-(2-fluorobenzyl)carbamoyl]-3,9-diazabicyclo[3.3.1]non-6-en-3-yl}-5-oxopentanoic acid methyl ester formate salt

20 Synthesized according to typical procedures A and E from bicyclononene AL23 and glutaric acid monomethylester chloride. LC-MS: R_t = 0.93; ES+: 750.39.

Example 295

25 (*rac.*)-5-{(1*R*^{*}, 5*S*^{*})-7-[4-[3-(2-Bromo-5-fluorophenoxy)propyl]phenyl]-6-[cyclopropyl-(3-trifluoromethylbenzyl)carbamoyl]-3,9-diazabicyclo[3.3.1]-non-6-en-3-yl}-5-oxopentanoic acid methyl ester formate salt

Synthesized according to typical procedures A and E from bicyclononene AL24 and glutaric acid monomethylester chloride. LC-MS: R_t = 0.96; ES+: 801.40.

30

Example 296

(rac.)-5-((1R, 5S*)-7-{4-[3-(2-Bromo-5-fluorophenoxy)propyl]phenyl}-6-[cyclopropyl-(2-methylbenzyl)carbamoyl]-3,9-diazabicyclo[3.3.1]non-6-en-3-yl}-5-oxopentanoic acid methyl ester formate salt*

5 Synthesized according to typical procedures A and E from bicyclononene AL25 and glutaric acid monomethylester chloride. LC-MS: R_t = 0.94; ES+: 746.43.

Example 297

10 *(rac.)-5-((1R*, 5S*)-7-{4-[3-(2-Bromo-5-fluorophenoxy)propyl]phenyl}-6-[cyclopropyl-[2-(4-methoxyphenoxy)ethyl]carbamoyl]-3,9-diazabicyclo[3.3.1]non-6-en-3-yl}-5-oxopentanoic acid methyl ester formate salt*

15 Synthesized according to typical procedures A and E from bicyclononene AL26 and glutaric acid monomethylester chloride. LC-MS: R_t = 0.93; ES+: 793.40.

Example 298

20 *(rac.)-5-((1R*, 5S*)-7-{4-[3-(2-Bromo-5-fluorophenoxy)propyl]phenyl}-6-[cyclopropyl-[2-(3-methylphenoxy)ethyl]carbamoyl]-3,9-diazabicyclo[3.3.1]non-6-en-3-yl}-5-oxopentanoic acid methyl ester formate salt*

25 Synthesized according to typical procedures A and E from bicyclononene AL27 and glutaric acid monomethylester chloride. LC-MS: R_t = 0.95; ES+: 776.41.

Example 299

30 *(rac.)-5-((1R*, 5S*)-7-{4-[3-(2-Bromo-5-fluorophenoxy)propyl]phenyl}-6-[cyclopropyl-[2-(3,4-dimethylphenoxy)ethyl]carbamoyl]-3,9-diazabicyclo[3.3.1]non-6-en-3-yl}-5-oxopentanoic acid methyl ester formate salt*

Synthesized according to typical procedures A and E from bicyclononene AL28 and glutaric acid monomethylester chloride. LC-MS: R_t = 0.97; ES+: 791.40.

Example 300

5

*(rac.)-5-[(1*R*^{*}, 5*S*^{*})-7-{4-[3-(2-Bromo-5-fluorophenoxy)propyl]phenyl}-6-(cyclopropylphenethylcarbamoyl)-3,9-diazabicyclo[3.3.1]non-6-en-3-yl]-5-oxo-pentanoic acid methyl ester formate salt*

10 Synthesized according to typical procedures A and E from bicyclononene AL29 and glutaric acid monomethylester chloride. LC-MS: R_t = 0.93; ES+: 746.43.

Example 301

15 *(rac.)-5-((1*R*^{*}, 5*S*^{*})-7-{4-[3-(2-Bromo-5-fluorophenoxy)propyl]phenyl}-6-{[2-(2-chlorophenyl)ethyl]cyclopropylcarbamoyl}-3,9-diazabicyclo[3.3.1]non-6-en-3-yl)-5-oxopentanoic acid methyl ester formate salt*

20 Synthesized according to typical procedures A and E from bicyclononene AL30 and glutaric acid monomethylester chloride. LC-MS: R_t = 0.95; ES+: 780.38.

Example 302

25 *(rac.)-5-((1*R*^{*}, 5*S*^{*})-7-{4-[3-(2-Bromo-5-fluorophenoxy)propyl]phenyl}-6-{[2-(2,3-difluorophenyl)ethyl]cyclopropylcarbamoyl}-3,9-diazabicyclo[3.3.1]non-6-en-3-yl)-5-oxopentanoic acid methyl ester formate salt*

30 Synthesized according to typical procedures A and E from bicyclononene AL31 and glutaric acid monomethylester chloride. LC-MS: R_t = 0.94; ES+: 782.40.

Example 303

(*rac.*)-5-((*1R*^{*}, *5S*^{*})-7-{4-[3-(2-Bromo-5-fluorophenoxy)propyl]phenyl}-6-{[2-(4-fluorophenyl)ethyl]cyclopropylcarbamoyl}-3,9-diazabicyclo[3.3.1]non-6-en-3-yl)-5-oxopentanoic acid methyl ester formate salt

- 5 Synthesized according to typical procedures A and E from bicyclononene **AL32** and glutaric acid monomethylester chloride. LC-MS: R_t = 0.94; ES+: 764.41.

Example 304

- 10 **(*rac.*)-5-((*1R*^{*}, *5S*^{*})-7-{4-[3-(2-Bromo-5-fluorophenoxy)propyl]phenyl}-6-{[2-(2-methylphenyl)ethyl]cyclopropylcarbamoyl}-3,9-diazabicyclo[3.3.1]non-6-en-3-yl)-5-oxopentanoic acid methyl ester**

Synthesized according to typical procedures A and E from bicyclononene **AL33**
15 and glutaric acid monomethylester chloride. LC-MS: R_t = 0.95; ES+: 760.43.

Example 305

- 20 **(*rac.*)-5-{(*1R*^{*}, *5S*^{*})-7-{4-[3-(2-Bromo-5-fluorophenoxy)propyl]phenyl}-6-[cyclopropyl-(3,5-dimethoxybenzyl)carbamoyl]-3,9-diazabicyclo[3.3.1]non-6-en-3-yl}-5-oxopentanoic acid methyl ester formate salt**

Synthesized according to typical procedures A and E from bicyclononene **AL35**
and glutaric acid monomethylester chloride. LC-MS: R_t = 0.93; ES+: 793.40.

25

Example 306

- 30 **(*rac.*)-5-((*1R*^{*}, *5S*^{*})-7-{4-[3-(2-Bromo-5-fluorophenoxy)propyl]phenyl}-6-{[2-(4-methylphenyl)ethyl]cyclopropylcarbamoyl}-3,9-diazabicyclo[3.3.1]non-6-en-3-yl)-5-oxopentanoic acid methyl ester formate salt**

Synthesized according to typical procedures A and E from bicyclononene AL36 and glutaric acid monomethylester chloride. LC-MS: R_t = 0.95; ES+: 760.42.

- Example 307

5

(*rac.*)-5-{(1*R*^{*}, 5*S*^{*})-7-{4-[3-(2-Bromo-5-fluorophenoxy)propyl]phenyl}-6-[(2-chlorobenzyl)ethylcarbamoyl]-3,9-diazabicyclo[3.3.1]non-6-en-3-yl}-2,2-dimethyl-5-oxopentanoic acid formate salt

10 Synthesized according to typical procedures K and E from bicyclononene AL22 and 2,2-dimethylglutaric anhydride. LC-MS: R_t = 0.93; ES+: 768.36.

Example 308

15 (*rac.*)-5-{(1*R*^{*}, 5*S*^{*})-7-{4-[3-(2-Bromo-5-fluorophenoxy)propyl]phenyl}-6-[cyclopropyl-(2-fluorobenzyl)carbamoyl]-3,9-diazabicyclo[3.3.1]non-6-en-3-yl}-2,2-dimethyl-5-oxo-pentanoic acid formate salt

20 Synthesized according to typical procedures K and E from bicyclononene AL23 and 2,2-dimethylglutaric anhydride. LC-MS: R_t = 0.92; ES+: 764.39.

Example 309

25 (*rac.*)-5-{(1*R*^{*}, 5*S*^{*})-7-{4-[3-(2-Bromo-5-fluorophenoxy)propyl]phenyl}-6-[cyclopropyl-(3-trifluoromethylbenzyl)carbamoyl]-3,9-diazabicyclo[3.3.1]non-6-en-3-yl}-2,2-dimethyl-5-oxopentanoic acid formate salt

Synthesized according to typical procedures K and E from bicyclononene AL24 and 2,2-dimethylglutaric anhydride. LC-MS: R_t = 0.94; ES+: 815.40.

30

Example 310

(rac.)-5-((1R, 5S*)-7-{4-[3-(2-Bromo-5-fluorophenoxy)propyl]phenyl}-6-[cyclopropyl-(2-methylbenzyl)carbamoyl]-3,9-diazabicyclo[3.3.1]non-6-en-3-yl)-2,2-dimethyl-5-oxo-pentanoic acid formate salt*

- 5 Synthesized according to typical procedures K and E from bicyclononene **AL25** and 2,2-dimethylglutaric anhydride. LC-MS: R_t = 0.93; ES+: 760.43.

Example 311

10 *(rac.)-5-((1R*, 5S*)-7-{4-[3-(2-Bromo-5-fluorophenoxy)propyl]phenyl}-6-[cyclopropyl-[2-(4-methoxyphenoxy)ethyl]-carbamoyl]-3,9-diazabicyclo-[3.3.1]non-6-en-3-yl)-2,2-dimethyl-5-oxopentanoic acid formate salt*

15 Synthesized according to typical procedures K and E from bicyclononene **AL26** and 2,2-dimethylglutaric anhydride. LC-MS: R_t = 0.92; ES+: >805.

Example 312

20 *(rac.)-5-((1R*, 5S*)-7-{4-[3-(2-Bromo-5-fluorophenoxy)propyl]phenyl}-6-[cyclopropyl-[2-(3-methylphenoxy)ethyl]-carbamoyl]-3,9-diazabicyclo[3.3.1]-non-6-en-3-yl)-2,2-dimethyl-5-oxopentanoic acid formate salt*

25 Synthesized according to typical procedures K and E from bicyclononene **AL27** and 2,2-dimethylglutaric anhydride. LC-MS: R_t = 0.94; ES+: 790.47.

Example 313

30 *(rac.)-5-((1R*, 5S*)-7-{4-[3-(2-Bromo-5-fluorophenoxy)propyl]phenyl}-6-[cyclopropyl-[2-(3,4-dimethylphenoxy)ethyl]-carbamoyl]-3,9-diazabicyclo-[3.3.1]non-6-en-3-yl)-2,2-dimethyl-5-oxopentanoic acid formate salt*

Synthesized according to typical procedures K and E from bicyclononene **AL28** and 2,2-dimethylglutaric anhydride. LC-MS: R_t = 0.95; ES+: 805.4.

Example 314

5

(*rac.*)-5-[(*1R*^{*}, *5S*^{*})-7-{4-[3-(2-Bromo-5-fluorophenoxy)propyl]phenyl}-6-(cyclopropylphenethylcarbamoyl)-3,9-diazabicyclo[3.3.1]non-6-en-3-yl]-2,2-dimethyl-5-oxopentanoic acid formate salt

10 Synthesized according to typical procedures K and E from bicyclononene **AL29** and 2,2-dimethylglutaric anhydride. LC-MS: R_t = 0.92; ES+: 760.43.

Example 315

15 (*rac.*)-5-((*1R*^{*}, *5S*^{*})-7-{4-[3-(2-Bromo-5-fluorophenoxy)propyl]phenyl}-6-{{[2-(2-chlorophenyl)ethyl]cyclopropylcarbamoyl}-3,9-diazabicyclo[3.3.1]non-6-en-3-yl}-2,2-dimethyl-5-oxopentanoic acid formate salt

20 Synthesized according to typical procedures K and E from bicyclononene **AL30** and 2,2-dimethylglutaric anhydride. LC-MS: R_t = 0.94; ES+: 794.41.

Example 316

25 (*rac.*)-5-((*1R*^{*}, *5S*^{*})-7-{4-[3-(2-Bromo-5-fluorophenoxy)propyl]phenyl}-6-{{[2-(2,3-difluorophenyl)ethyl]cyclopropylcarbamoyl}-3,9-diazabicyclo[3.3.1]non-6-en-3-yl}-2,2-dimethyl-5-oxopentanoic acid formate salt

30 Synthesized according to typical procedures K and E from bicyclononene **AL31** and 2,2-dimethylglutaric anhydride. LC-MS: R_t = 0.93; ES+: 796.44.

Example 317

(rac.)-5-((1R, 5S*)-7-{4-[3-(2-Bromo-5-fluorophenoxy)propyl]phenyl}-6-[(2-(4-fluorophenyl)ethyl]cyclopropylcarbamoyl}-3,9-diazabicyclo[3.3.1]non-6-en-3-yl)-2,2-dimethyl-5-oxopentanoic acid formate salt*

- 5 Synthesized according to typical procedures K and E from bicyclononene AL32 and 2,2-dimethylglutaric anhydride. LC-MS: R_t = 0.93; ES+: 778.42.

Example 318

- 10 *(rac.)-5-((1R*, 5S*)-7-{4-[3-(2-Bromo-5-fluorophenoxy)propyl]phenyl}-6-[(2-methylphenyl)ethyl]cyclopropylcarbamoyl}-3,9-diazabicyclo[3.3.1]non-6-en-3-yl)-2,2-dimethyl-5-oxopentanoic acid formate salt*

Synthesized according to typical procedures K and E from bicyclononene AL33
15 and 2,2-dimethylglutaric anhydride. LC-MS: R_t = 0.94; ES+: 774.46.

Example 319

- 20 *(rac.)-5-((1R*, 5S*)-7-{4-[3-(2-Bromo-5-fluorophenoxy)propyl]phenyl}-6-[cyclopropyl-(3,5-dimethoxybenzyl)carbamoyl]-3,9-diazabicyclo[3.3.1]non-6-en-3-yl)-2,2-dimethyl-5-oxopentanoic acid formate salt*

Synthesized according to typical procedures K and E from bicyclononene AL35
25 and 2,2-dimethylglutaric anhydride. LC-MS: R_t = 0.92; ES+: >805.

Example 320

- 30 *(rac.)-5-((1R*, 5S*)-7-{4-[3-(2-Bromo-5-fluorophenoxy)propyl]phenyl}-6-[(2-(4-methylphenyl)ethyl]cyclopropylcarbamoyl}-3,9-diazabicyclo[3.3.1]non-6-en-3-yl)-2,2-dimethyl-5-oxopentanoic acid formate salt*

Synthesized according to typical procedures K and E from bicyclononene AL36 and 2,2-dimethylglutaric anhydride. LC-MS: R_t = 0.94; ES+: 774.44.

Example 321

5

1:1-Mixture of (rac.)-(2*R*^{*}, 3*S*^{*})-4-{(1*R*^{*}, 5*S*^{*})-7-[4-[3-(2-bromo-5-fluoro-phenoxy)propyl]phenyl]-6-[(2-chlorobenzyl)ethylcarbamoyl]-3,9-diazabi-cyclo[3.3.1]non-6-en-3-yl}-2,3-dihydroxy-4-oxobutyric acid formate salt and (rac.)-(2*S*^{*}, 3*R*^{*})-4-{(1*R*^{*}, 5*S*^{*})-7-[4-[3-(2-bromo-5-fluorophenoxy)propyl]-phenyl]-6-[(2-chlorobenzyl)ethylcarbamoyl]-3,9-diazabicyclo-[3.3.1]non-6-en-3-yl}-2,3-dihydroxy-4-oxobutyric acid formate salt

10

Synthesized according to typical procedures K and E from bicyclononene AL22 and (S,R)-2,3-dihydroxysuccinic anhydride. LC-MS: R_t = 0.87; ES+: 758.33.

15

Example 322

1:1-Mixture of (rac.)-(2*R*^{*}, 3*S*^{*})-4-{(1*R*^{*}, 5*S*^{*})-7-[4-[3-(2-bromo-5-fluoro-phenoxy)propyl]phenyl]-6-[cyclopropyl-(2-fluorobenzyl)carbamoyl]-3,9-diazabicyclo[3.3.1]non-6-en-3-yl}-2,3-dihydroxy-4-oxobutyric acid formate salt and (rac.)-(2*S*^{*}, 3*R*^{*})-4-{(1*R*^{*}, 5*S*^{*})-7-[4-[3-(2-bromo-5-fluoro-phenoxy)propyl]phenyl]-6-[cyclopropyl-(2-fluorobenzyl)carbamoyl]-3,9-diazabicyclo[3.3.1]non-6-en-3-yl}-2,3-dihydroxy-4-oxobutyric acid formate salt

25

Synthesized according to typical procedures K and E from bicyclononene AL23 and (S,R)-2,3-dihydroxysuccinic anhydride. LC-MS: R_t = 0.85; ES+: 754.37.

30

1:1-Mixture of (rac.)-(2*R*^{*}, 3*S*^{*})-4-{(1*R*^{*}, 5*S*^{*})-7-[4-[3-(2-bromo-5-fluoro-phenoxy)propyl]phenyl]-6-[cyclopropyl-(3-trifluoromethylbenzyl)-

Example 323

carbamoyl]-3,9-diazabicyclo[3.3.1]non-6-en-3-yl}-2,3-dihydroxy-4-oxobutyric acid formate salt and (*rac.*)-(2*S*^{*,} 3*R*^{*)}-4-{(1*R*^{*,} 5*S*^{*)}-7-{4-[3-(2-bromo-5-fluorophenoxy)propyl]-phenyl}-6-[cyclopropyl-(3-trifluoromethylbenzyl)-carbamoyl]-3,9-diazabicyclo[3.3.1]non-6-en-3-yl}-2,3-dihydroxy-4-oxobutyric acid formate salt

Synthesized according to typical procedures K and E from bicyclononene AL24 and (S,R)-2,3-dihydroxysuccinic anhydride. LC-MS: R_t = 0.89; ES+: 805.4.

10 **Example 324**

1:1-Mixture of (*rac.*)-(2*R*^{*,} 3*S*^{*)}-4-{(1*R*^{*,} 5*S*^{*)}-7-{4-[3-(2-bromo-5-fluorophenoxy)propyl]phenyl}-6-[cyclopropyl-(2-methylbenzyl)carbamoyl]-3,9-diazabicyclo[3.3.1]non-6-en-3-yl}-2,3-dihydroxy-4-oxobutyric acid formate salt and (*rac.*)-(2*S*^{*,} 3*R*^{*)}-4-{(1*R*^{*,} 5*S*^{*)}-7-{4-[3-(2-bromo-5-fluorophenoxy)propyl]phenyl}-6-[cyclopropyl-(2-methylbenzyl)carbamoyl]-3,9-diazabicyclo[3.3.1]non-6-en-3-yl}-2,3-dihydroxy-4-oxobutyric acid formate salt

Synthesized according to typical procedures K and E from bicyclononene AL25 and (S,R)-2,3-dihydroxysuccinic anhydride. LC-MS: R_t = 0.87; ES+: 750.39.

Example 325

1:1-Mixture of (*rac.*)-(2*R*^{*,} 3*S*^{*)}-4-{(1*R*^{*,} 5*S*^{*)}-7-{4-[3-(2-bromo-5-fluorophenoxy)propyl]phenyl}-6-[cyclopropyl-[2-(4-methoxyphenoxy)ethyl]-carbamoyl]-3,9-diazabicyclo[3.3.1]non-6-en-3-yl}-2,3-dihydroxy-4-oxobutyric acid formate salt and (*rac.*)-(2*S*^{*,} 3*R*^{*)}-4-{(1*R*^{*,} 5*S*^{*)}-7-{4-[3-(2-bromo-5-fluorophenoxy)propyl]phenyl}-6-[cyclopropyl-[2-(4-methoxyphenoxy)ethyl]-carbamoyl]-3,9-diazabicyclo[3.3.1]non-6-en-3-yl}-2,3-dihydroxy-4-oxobutyric acid formate salt

Synthesized according to typical procedures K and E from bicyclononene **AL26** and (S,R)-2,3-dihydroxysuccinic anhydride. LC-MS: R_t = 0.86; ES+: 796.42.

Example 326

5

1:1-Mixture of (*rac.*)-(2*R*^{*,} 3*S*^{*)}-4-((1*R*^{*,} 5*S*^{*)}-7-{4-[3-(2-bromo-5-fluoro-phenoxy)propyl]phenyl}-6-{cyclopropyl-[2-(3-methylphenoxy)ethyl]-carbamoyl}-3,9-diazabicyclo[3.3.1]non-6-en-3-yl)-2,3-dihydroxy-4-oxobutyric acid formate salt and (*rac.*)-(2*S*^{*,} 3*R*^{*)}-4-((1*R*^{*,} 5*S*^{*)}-7-{4-[3-(2-bromo-5-fluorophenoxy)propyl]phenyl}-6-{cyclopropyl-[2-(3-methylphenoxy)ethyl]-carbamoyl}-3,9-diazabicyclo[3.3.1]non-6-en-3-yl)-2,3-dihydroxy-4-oxobutyric acid formate salt

10

Synthesized according to typical procedures K and E from bicyclononene **AL27** and (S,R)-2,3-dihydroxysuccinic anhydride. LC-MS: R_t = 0.88; ES+: 780.38.

15

Example 327

20

1:1-Mixture of (*rac.*)-(2*R*^{*,} 3*S*^{*)}-4-((1*R*^{*,} 5*S*^{*)}-7-{4-[3-(2-bromo-5-fluoro-phenoxy)propyl]phenyl}-6-{cyclopropyl-[2-(3,4-dimethylphenoxy)ethyl]-carbamoyl}-3,9-diazabicyclo[3.3.1]non-6-en-3-yl)-2,3-dihydroxy-4-oxobutyric acid formate salt and (*rac.*)-(2*S*^{*,} 3*R*^{*)}-4-((1*R*^{*,} 5*S*^{*)}-7-{4-[3-(2-bromo-5-fluorophenoxy)-propyl]phenyl}-6-{cyclopropyl-[2-(3,4-dimethylphenoxy)-ethyl]carbamoyl}-3,9-diazabicyclo[3.3.1]non-6-en-3-yl)-2,3-dihydroxy-4-oxobutyric acid formate salt

25

Synthesized according to typical procedures K and E from bicyclononene **AL28** and (S,R)-2,3-dihydroxysuccinic anhydride. LC-MS: R_t = 0.89; ES+: 794.44.

30

Example 328

1:1-Mixture of (*rac.*)-(2*R*^{*}, 3*S*^{*})-4-[(*IR*^{*}, 5*S*^{*})-7-{4-[3-(2-bromo-5-fluoro-phenoxy)propyl]phenyl}-6-(cyclopropylphenethylcarbamoyl)-3,9-diazabi-cyclo[3.3.1]non-6-en-3-yl]-2,3-dihydroxy-4-oxobutyric acid formate salt and (*rac.*)-(2*S*^{*}, 3*R*^{*})-4-[(*IR*^{*}, 5*S*^{*})-7-{4-[3-(2-bromo-5-fluorophenoxy)propyl]-phenyl}-6-(cyclopropylphenethylcarbamoyl)-3,9-diazabicyclo[3.3.1]non-6-en-3-yl]-2,3-dihydroxy-4-oxobutyric acid formate salt

Synthesized according to typical procedures K and E from bicyclononene AL29 and (S,R)-2,3-dihydroxysuccinic anhydride. LC-MS: R_t = 0.86; ES+: 750.36.

10

Example 329

1:1-Mixture of (*rac.*)-(2*R*^{*}, 3*S*^{*})-4-[(*IR*^{*}, 5*S*^{*})-7-{4-[3-(2-bromo-5-fluoro-phenoxy)propyl]phenyl}-6-{[2-(2-chlorophenyl)ethyl]cyclopropylcarbamoyl}-3,9-diazabicyclo[3.3.1]non-6-en-3-yl]-2,3-dihydroxy-4-oxobutyric acid formate salt and (*rac.*)-(2*S*^{*}, 3*R*^{*})-4-[(*IR*^{*}, 5*S*^{*})-7-{4-[3-(2-bromo-5-fluorophenoxy)propyl]-phenyl}-6-{[2-(2-chlorophenyl)ethyl]cyclopropyl-carbamoyl}-3,9-diazabicyclo[3.3.1]non-6-en-3-yl]-2,3-dihydroxy-4-oxobutyric acid formate salt

20

Synthesized according to typical procedures K and E from bicyclononene AL30 and (S,R)-2,3-dihydroxysuccinic anhydride. LC-MS: R_t = 0.88; ES+: 784.34.

25

1:1-Mixture of (*rac.*)-(2*R*^{*}, 3*S*^{*})-4-[(*IR*^{*}, 5*S*^{*})-7-{4-[3-(2-bromo-5-fluoro-phenoxy)propyl]phenyl}-6-{[2-(4-fluorophenyl)ethyl]cyclopropylcarbamoyl}-3,9-diazabicyclo[3.3.1]non-6-en-3-yl]-2,3-dihydroxy-4-oxobutyric acid formate salt and (*rac.*)-(2*S*^{*}, 3*R*^{*})-4-[(*IR*^{*}, 5*S*^{*})-7-{4-[3-(2-bromo-5-fluorophenoxy)propyl]-phenyl}-6-{[2-(4-fluorophenyl)ethyl]cyclopropyl-carbamoyl}-3,9-diazabicyclo[3.3.1]non-6-en-3-yl]-2,3-dihydroxy-4-oxobutyric acid formate salt

Synthesized according to typical procedures K and E from bicyclononene AL32 and (S,R)-2,3-dihydroxysuccinic anhydride. LC-MS: R_t = 0.85; ES+: 768.34.

5 Example 331

1:1-Mixture of (*rac.*)-(2*R*^{*}, 3*S*^{*})-4-((1*R*^{*}, 5*S*^{*})-7-{4-[3-(2-bromo-5-fluoro-phenoxy)propyl]phenyl}-6-{[2-(2-methylphenyl)ethyl]cyclopropyl-carbamoyl}-3,9-diazabicyclo[3.3.1]non-6-en-3-yl)-2,3-dihydroxy-4-oxobutyric acid formate salt and (*rac.*)-(2*S*^{*}, 3*R*^{*})-4-((1*R*^{*}, 5*S*^{*})-7-{4-[3-(2-bromo-5-fluorophenoxy)propyl]phenyl}-6-{[2-(2-methylphenyl)ethyl]cyclopropyl-carbamoyl}-3,9-diazabicyclo-[3.3.1]non-6-en-3-yl)-2,3-dihydroxy-4-oxobutyric acid formate salt

15 Synthesized according to typical procedures K and E from bicyclononene AL33 and (S,R)-2,3-dihydroxysuccinic anhydride. LC-MS: R_t = 0.88; ES+: 764.41.

Example 332

20 1:1-Mixture of (*rac.*)-(2*R*^{*}, 3*S*^{*})-4-((1*R*^{*}, 5*S*^{*})-7-{4-[3-(2-bromo-5-fluoro-phenoxy)propyl]phenyl}-6-[cyclopropyl-(3,5-dimethoxybenzyl)carbamoyl]-3,9-diazabicyclo[3.3.1]non-6-en-3-yl)-2,3-dihydroxy-4-oxobutyric acid formate salt and (*rac.*)-(2*S*^{*}, 3*R*^{*})-4-((1*R*^{*}, 5*S*^{*})-7-{4-[3-(2-bromo-5-fluoro-phenoxy)propyl]phenyl}-6-[cyclopropyl-(3,5-dimethoxybenzyl)carbamoyl]-3,9-diazabicyclo[3.3.1]non-6-en-3-yl)-2,3-dihydroxy-4-oxobutyric acid formate salt

Synthesized according to typical procedures K and E from bicyclononene AL35 and (S,R)-2,3-dihydroxysuccinic anhydride. LC-MS: R_t = 0.86; ES+: 797.38.

1:1-Mixture of (*rac.*)-(2*R*^{*,} 3*S*^{*})-4-((1*R*^{*,} 5*S*^{*})-7-{4-[3-(2-bromo-5-fluoro-phenoxy)propyl]phenyl}-6-{[2-(4-methylphenyl)ethyl]cyclopropyl-carbamoyl}-3,9-diazabicyclo[3.3.1]non-6-en-3-yl)-2,3-dihydroxy-4-oxobutyric acid formate salt and (*rac.*)-(2*S*^{*,} 3*R*^{*})-4-((1*R*^{*,} 5*S*^{*})-7-{4-[3-(2-bromo-5-fluorophenoxy)propyl]phenyl}-6-{[2-(4-methylphenyl)ethyl]cyclopropyl-carbamoyl}-3,9-diazabicyclo-[3.3.1]non-6-en-3-yl)-2,3-dihydroxy-4-oxobutyric acid formate salt

10 Synthesized according to typical procedures K and E from bicyclononene AL36 and (S,R)-2,3-dihydroxysuccinic anhydride. LC-MS: R_t = 0.87; ES+: 764.38.

Example 334

15 (*rac.*)-(1*R*^{*,} 5*S*^{*})-7-{4-[3-(2-Bromo-5-fluorophenoxy)propyl]phenyl}-3-(4-carbamoylbutyryl)-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid (2-chlorobenzyl)ethylamide formate salt

20 Synthesized according to typical procedures G and E from bicyclononene AL22 and 4-carbamoylbutyric acid. LC-MS: R_t = 0.87; ES+: 739.49.

Example 335

25 (*rac.*)-(1*R*^{*,} 5*S*^{*})-7-{4-[3-(2-Bromo-5-fluorophenoxy)propyl]phenyl}-3-(4-carbamoylbutyryl)-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid cyclopropyl-(2-fluorobenzyl)amide formate salt

30 Synthesized according to typical procedures G and E from bicyclononene AL23 and 4-carbamoylbutyric acid. LC-MS: R_t = 0.86; ES+: 735.50.

Example 336

(*rac.*)-(1*R*^{*}, 5*S*^{*})-7-{4-[3-(2-Bromo-5-fluorophenoxy)propyl]phenyl}-3-(4-carbamoylbutyryl)-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid cyclopropyl-(2-methylbenzyl)amide formate salt

- 5 Synthesized according to typical procedures G and E from bicyclononene AL25 and 4-carbamoylbutyric acid. LC-MS: R_t = 0.88; ES+: 731.55.

Example 337

10 (*rac.*)-(1*R*^{*}, 5*S*^{*})-7-{4-[3-(2-Bromo-5-fluorophenoxy)propyl]phenyl}-3-(4-carbamoylbutyryl)-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid cyclopropyl-[2-(4-methoxyphenoxy)ethyl]amide formate salt

- 15 Synthesized according to typical procedures G and E from bicyclononene AL26 and 4-carbamoylbutyric acid. LC-MS: R_t = 0.86; ES+: 777.52.

Example 338

20 (*rac.*)-(1*R*^{*}, 5*S*^{*})-7-{4-[3-(2-Bromo-5-fluorophenoxy)propyl]phenyl}-3-(4-carbamoylbutyryl)-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid cyclopropyl-[2-(3,4-dimethylphenoxy)ethyl]amide formate salt

- 25 Synthesized according to typical procedures G and E from bicyclononene AL28 and 4-carbamoylbutyric acid. LC-MS: R_t = 0.90; ES+: 775.55.

25

Example 339

30 (*rac.*)-(1*R*^{*}, 5*S*^{*})-7-{4-[3-(2-Bromo-5-fluorophenoxy)propyl]phenyl}-3-(4-carbamoylbutyryl)-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid [2-(2-chlorophenyl)ethyl]cyclopropylamide formate salt

Synthesized according to typical procedures G and E from bicyclononene AL30 and 4-carbamoylbutyric acid. LC-MS: R_t = 0.89; ES+: 765.46.

Example 340

5

(*rac.*)-(IR*, 5S*)-7-{4-[3-(2-Bromo-5-fluorophenoxy)propyl]phenyl}-3-(4-carbamoylbutyryl)-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid [2-(2,3-difluorophenyl)ethyl]cyclopropylamide formate salt

10 Synthesized according to typical procedures G and E from bicyclononene AL31 and 4-carbamoylbutyric acid. LC-MS: R_t = 0.88; ES+: 767.48.

Example 341

15 (*rac.*)-(IR*, 5S*)-7-{4-[3-(2-Bromo-5-fluorophenoxy)propyl]phenyl}-3-(4-carbamoylbutyryl)-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid [2-(4-fluorophenyl)ethyl]cyclopropylamide formate salt

20 Synthesized according to typical procedures G and E from bicyclononene AL32 and 4-carbamoylbutyric acid. LC-MS: R_t = 0.87; ES+: 749.52.

Example 342

25 (*rac.*)-(IR*, 5S*)-7-{4-[3-(2-Bromo-5-fluorophenoxy)propyl]phenyl}-3-(4-carbamoylbutyryl)-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid [2-(2-methylphenyl)ethyl]cyclopropylamide formate salt

Synthesized according to typical procedures G and E from bicyclononene AL33 and 4-carbamoylbutyric acid. LC-MS: R_t = 0.89; ES+: 745.54.

30

Example 343

(rac.)-(1R, 5S*)-7-{4-[3-(2-Bromo-5-fluorophenoxy)propyl]phenyl}-3-(4-carbamoylbutyryl)-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid cyclopropyl-(3,5-dimethoxybenzyl)amide formate salt*

5 Synthesized according to typical procedures G and E from bicyclononene AL35 and 4-carbamoylbutyric acid. LC-MS: R_t = 0.86; ES+: 777.53.

Example 344

10 *(rac.)-(1R*, 5S*)-7-{4-[3-(2-Bromo-5-fluorophenoxy)propyl]phenyl}-3-(4-carbamoylbutyryl)-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid [2-(4-methylphenyl)ethyl]cyclopropylamide formate salt*

15 Synthesized according to typical procedures G and E from bicyclononene AL36 and 4-carbamoylbutyric acid. LC-MS: R_t = 0.88; ES+: 745.54.

Example 345

20 *1:1-Mixture of (rac.)-(3R*)-5-{(1R*, 5S*)-7-{4-[3-(2-bromo-5-fluorophenoxy)-propyl]phenyl}-6-[(2-chlorobenzyl)ethylcarbamoyl]-3,9-diazabicyclo[3.3.1]-non-6-en-3-yl}-3-hydroxy-5-oxopentanoic acid formate salt and (rac.)-(3S*)-5-{(1R*, 5S*)-7-{4-[3-(2-bromo-5-fluorophenoxy)-propyl]phenyl}-6-[(2-chlorobenzyl)ethylcarbamoyl]-3,9-diazabicyclo[3.3.1]non-6-en-3-yl}-3-hydroxy-5-oxopentanoic acid formate salt*

25 Synthesized according to typical procedures K, E and L from bicyclononene AL22 and 3-(*tert*-butyldimethylsilyloxy)glutaric anhydride. LC-MS: R_t = 0.87; ES+: 756.44.

30 **Example 346**

1:1-Mixture of (rac.)-(3R*)-5-{(1R*, 5S*)-7-{4-[3-(2-bromo-5-fluorophenoxy)-propyl]phenyl}-6-[cyclopropyl-(2-fluorobenzyl)carbamoyl]-3,9-diazabicyclo[3.3.1]non-6-en-3-yl}-3-hydroxy-5-oxopentanoic acid formate salt and (rac.)-(3S*)-5-{(1R*, 5S*)-7-{4-[3-(2-bromo-5-fluorophenoxy)propyl]phenyl}-6-[cyclopropyl-(2-fluorobenzyl)carbamoyl]-3,9-diazabicyclo[3.3.1]non-6-en-3-yl}-3-hydroxy-5-oxopentanoic acid formate salt

Synthesized according to typical procedures K, E and L from bicyclononene AL23 and 3-(*tert*-butyldimethylsilyloxy)glutaric anhydride. LC-MS: R_t = 0.86; ES+: 752.46.

Example 347

1:1-Mixture of (rac.)-(3R*)-5-{(1R*, 5S*)-7-{4-[3-(2-bromo-5-fluorophenoxy)-propyl]phenyl}-6-[cyclopropyl-(3-trifluoromethylbenzyl)carbamoyl]-3,9-diazabicyclo[3.3.1]non-6-en-3-yl}-3-hydroxy-5-oxopentanoic acid formate salt and (rac.)-(3S*)-5-{(1R*, 5S*)-7-{4-[3-(2-bromo-5-fluorophenoxy)propyl]phenyl}-6-[cyclopropyl-(3-trifluoromethylbenzyl)carbamoyl]-3,9-diazabicyclo[3.3.1]non-6-en-3-yl}-3-hydroxy-5-oxopentanoic acid formate salt

Synthesized according to typical procedures K, E and L from bicyclononene AL24 and 3-(*tert*-butyldimethylsilyloxy)glutaric anhydride. LC-MS: R_t = 0.89; ES+: 803.40.

Example 348

1:1-Mixture of (rac.)-(3R*)-5-{(1R*, 5S*)-7-{4-[3-(2-bromo-5-fluorophenoxy)-propyl]phenyl}-6-[cyclopropyl-(2-methylbenzyl)carbamoyl]-3,9-diazabicyclo[3.3.1]non-6-en-3-yl}-3-hydroxy-5-oxopentanoic acid formate salt and (rac.)-(3S*)-5-{(1R*, 5S*)-7-{4-[3-(2-bromo-5-fluorophenoxy)propyl]phenyl}-6-[cyclopropyl-(2-methylbenzyl)carbamoyl]-3,9-diazabicyclo[3.3.1]non-6-en-3-yl}-3-hydroxy-5-oxopentanoic acid formate salt

Synthesized according to typical procedures K, E and L from bicyclononene AL25 and 3-(*tert*-butyldimethylsilyloxy)glutaric anhydride. LC-MS: R_t = 0.87; ES+: 748.52.

5

Example 349

- 1:1-Mixture of (*rac.*)-(3*R*^{*})-5-((1*R*^{*}, 5*S*^{*})-7-{4-[3-(2-bromo-5-fluorophenoxy)propyl]phenyl}-6-{cyclopropyl-[2-(4-methoxyphenoxy)ethyl]carbamoyl}-3,9-diazabicyclo[3.3.1]non-6-en-3-yl)-3-hydroxy-5-oxopentanoic acid formate salt and (*rac.*)-(3*S*^{*})-5-((1*R*^{*}, 5*S*^{*})-7-{4-[3-(2-bromo-5-fluorophenoxy)propyl]phenyl}-6-{cyclopropyl-[2-(4-methoxyphenoxy)ethyl]carbamoyl}-3,9-diazabicyclo[3.3.1]non-6-en-3-yl)-3-hydroxy-5-oxopentanoic acid formate salt
- 15 Synthesized according to typical procedures K, E and L from bicyclononene AL26 and 3-(*tert*-butyldimethylsilyloxy)glutaric anhydride. LC-MS: R_t = 0.86; ES+: 794.40.

20

Example 350

- 1:1-Mixture of (*rac.*)-(3*R*^{*})-5-((1*R*^{*}, 5*S*^{*})-7-{4-[3-(2-bromo-5-fluorophenoxy)propyl]phenyl}-6-{cyclopropyl-[2-(3-methylphenoxy)ethyl]carbamoyl}-3,9-diazabicyclo[3.3.1]non-6-en-3-yl)-3-hydroxy-5-oxopentanoic acid formate salt and (*rac.*)-(3*S*^{*})-5-((1*R*^{*}, 5*S*^{*})-7-{4-[3-(2-bromo-5-fluorophenoxy)propyl]phenyl}-6-{cyclopropyl-[2-(3-methylphenoxy)ethyl]carbamoyl}-3,9-diazabicyclo[3.3.1]non-6-en-3-yl)-3-hydroxy-5-oxopentanoic acid formate salt

25

- Synthesized according to typical procedures K, E and L from bicyclononene AL27 and 3-(*tert*-butyldimethylsilyloxy)glutaric anhydride. LC-MS: R_t = 0.89; ES+: 778.52.

Example 351

1:1-Mixture of (*rac.*)-(3*R**)-5-((*IR**, 5*S**)-7-{4-[3-(2-bromo-5-fluorophenoxy)-propyl]phenyl}-6-{cyclopropyl-[2-(3,4-dimethylphenoxy)ethyl]carbamoyl}-3,9-diazabicyclo[3.3.1]non-6-en-3-yl)-3-hydroxy-5-oxopentanoic acid formate salt and (*rac.*)-(3*S**)-5-((*IR**, 5*S**)-7-{4-[3-(2-bromo-5-fluorophenoxy)propyl]-phenyl}-6-{cyclopropyl-[2-(3,4-dimethylphenoxy)ethyl]carbamoyl}-3,9-diazabicyclo[3.3.1]-non-6-en-3-yl)-3-hydroxy-5-oxopentanoic acid formate salt

Synthesized according to typical procedures K, E and L from bicyclononene AL28 and 3-(*tert*-butyldimethylsilyloxy)glutaric anhydride. LC-MS: R_t = 0.90; ES+: 792.42.

Example 352

1:1-Mixture of (*rac.*)-(3*R**)-5-[(*IR**, 5*S**)-7-{4-[3-(2-bromo-5-fluorophenoxy)-propyl]phenyl}-6-(cyclopropylphenethylcarbamoyl)-3,9-diazabicyclo[3.3.1]-non-6-en-3-yl]-3-hydroxy-5-oxopentanoic acid formate salt and (*rac.*)-(3*R**)-5-[(*IR**, 5*S**)-7-{4-[3-(2-bromo-5-fluorophenoxy)propyl]phenyl}-6-(cyclopropylphenethylcarbamoyl)-3,9-diazabicyclo[3.3.1]non-6-en-3-yl]-3-hydroxy-5-oxopentanoic acid formate salt

Synthesized according to typical procedures K, E and L from bicyclononene AL29 and 3-(*tert*-butyldimethylsilyloxy)glutaric anhydride. LC-MS: R_t = 0.86; ES+: 748.50.

25

Example 353

1:1-Mixture of (*rac.*)-(3*R**)-5-((*IR**, 5*S**)-7-{4-[3-(2-bromo-5-fluorophenoxy)-propyl]phenyl}-6-{[2-(2-chlorophenyl)ethyl]cyclopropylcarbamoyl}-3,9-diazabicyclo[3.3.1]non-6-en-3-yl)-3-hydroxy-5-oxopentanoic acid formate salt and (*rac.*)-(3*R**)-5-((*IR**, 5*S**)-7-{4-[3-(2-bromo-5-fluorophenoxy)propyl]-

phenyl}-6-{[2-(2-chlorophenyl)ethyl]cyclopropylcarbamoyl}-3,9-diaza-bicyclo[3.3.1]non-6-en-3-yl)-3-hydroxy-5-oxopentanoic acid formate salt

Synthesized according to typical procedures K, E and L from bicyclononene
5 AL30 and 3-(*tert*-butyldimethylsilyloxy)glutaric anhydride. LC-MS: R_t = 0.88;
ES+: 782.49.

Example 354

- 10 **1:1-Mixture of (*rac.*)-(3*R**)-5-((1*R**, 5*S**)-7-{4-[3-(2-bromo-5-fluorophenoxy)-propyl]phenyl}-6-{[2-(2,3-difluorophenyl)ethyl]cyclopropylcarbamoyl}-3,9-diazabicyclo[3.3.1]non-6-en-3-yl)-3-hydroxy-5-oxopentanoic acid formate salt and (*rac.*)-(3*R**)-5-((1*R**, 5*S**)-7-{4-[3-(2-bromo-5-fluorophenoxy)propyl]-phenyl}-6-{[2-(2,3-difluorophenyl)ethyl]cyclopropylcarbamoyl}-3,9-diazabi-15 cyclo[3.3.1]non-6-en-3-yl)-3-hydroxy-5-oxopentanoic acid formate salt**

Synthesized according to typical procedures K, E and L from bicyclononene AL31 and 3-(*tert*-butyldimethylsilyloxy)glutaric anhydride. LC-MS: R_t = 0.88;
ES+: 784.53.

20

Example 355

- 1:1-Mixture of (*rac.*)-(3*R**)-5-((1*R**, 5*S**)-7-{4-[3-(2-bromo-5-fluorophenoxy)-propyl]phenyl}-6-{[2-(4-fluorophenyl)ethyl]cyclopropylcarbamoyl}-3,9-diazabicyclo[3.3.1]non-6-en-3-yl)-3-hydroxy-5-oxopentanoic acid formate salt and (*rac.*)-(3*R**)-5-((1*R**, 5*S**)-7-{4-[3-(2-bromo-5-fluorophenoxy)propyl]phenyl}-6-{[2-(4-fluorophenyl)ethyl]cyclopropylcarbamoyl}-3,9-diazabicyclo[3.3.1]-25 non-6-en-3-yl)-3-hydroxy-5-oxopentanoic acid formate salt

- 30 Synthesized according to typical procedures K, E and L from bicyclononene AL32 and 3-(*tert*-butyldimethylsilyloxy)glutaric anhydride. LC-MS: R_t = 0.87;
ES+: 766.47.

Example 356

1:1-Mixture of (*rac.*)-(3*R**)-5-((*IR**, *S**)-7-{4-[3-(2-bromo-5-fluorophenoxy)-
5 propyl]phenyl}-6-{[2-(2-methylphenyl)ethyl]cyclopropylcarbamoyl}-3,9-diazabicyclo[3.3.1]non-6-en-3-yl)-3-hydroxy-5-oxopentanoic acid formate salt
and (*rac.*)-(3*R**)-5-((*IR**, *S**)-7-{4-[3-(2-bromo-5-fluorophenoxy)propyl]-
phenyl}-6-{[2-(2-methylphenyl)ethyl]cyclopropylcarbamoyl}-3,9-diazabi-
cyclo[3.3.1]non-6-en-3-yl)-3-hydroxy-5-oxopentanoic acid formate salt

10

Synthesized according to typical procedures K, E and L from bicyclononene AL33 and 3-(*tert*-butyldimethylsilyloxy)glutaric anhydride. LC-MS: R_t = 0.88;
ES+: 762.55.

Example 357

1:1-Mixture of (*rac.*)-(3*R**)-5-((*IR**, *S**)-7-{4-[3-(2-bromo-5-fluorophenoxy)-
propyl]phenyl}-6-[cyclopropyl-(3,5-dimethoxybenzyl)carbamoyl]-3,9-diaza-
bicyclo[3.3.1]non-6-en-3-yl)-3-hydroxy-5-oxopentanoic acid formate salt and
20 (*rac.*)-(3*S**)-5-((*IR**, *S**)-7-{4-[3-(2-bromo-5-fluorophenoxy)propyl]phenyl}-
6-[cyclopropyl-(3,5-dimethoxybenzyl)carbamoyl]-3,9-diazabicyclo-[3.3.1]non-
6-en-3-yl)-3-hydroxy-5-oxopentanoic acid formate salt

Synthesized according to typical procedures K, E and L from bicyclononene
25 AL35 and 3-(*tert*-butyldimethylsilyloxy)glutaric anhydride. LC-MS: R_t = 0.86;
ES+: 794.38.

Example 358

30 1:1-Mixture of (*rac.*)-(3*R**)-5-((*IR**, *S**)-7-{4-[3-(2-bromo-5-fluorophenoxy)-
propyl]phenyl}-6-{[2-(4-methylphenyl)ethyl]cyclopropylcarbamoyl}-3,9-
diaza-
bicyclo[3.3.1]non-6-en-3-yl)-3-hydroxy-5-oxopentanoic acid formate salt

and (*rac.*)-(3*R**)-5-((1*R**, 5*S**)-7-{4-[3-(2-bromo-5-fluorophenoxy)propyl]phenyl}-6-{[2-(4-methylphenyl)ethyl]cyclopropylcarbamoyl}-3,9-diazabicyclo[3.3.1]non-6-en-3-yl)-3-hydroxy-5-oxopentanoic acid formate salt

- 5 Synthesized according to typical procedures K, E and L from bicyclononene AL36 and 3-(*tert*-butyldimethylsilyloxy)glutaric anhydride. LC-MS: R_t = 0.88; ES+: 762.55.

Example 359

10

1:1-Mixture of (*rac.*)-(3*R**)-5-((1*R**, 5*S**)-6-{cyclopropyl-[2-(2-hydroxyethyl)benzyl]carbamoyl}-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-en-3-yl)-3-hydroxy-5-oxopentanoic acid formate salt and (*rac.*)-(3*S**)-5-((1*R**, 5*S**)-6-{cyclopropyl-[2-(2-hydroxyethyl)benzyl]carbamoyl}-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-en-3-yl)-3-hydroxy-5-oxopentanoic acid formate salt

Synthesized according to typical procedures K, E and L from bicyclononene AL19 and 3-(*tert*-butyldimethylsilyloxy)glutaric anhydride. LC-MS: R_t = 0.81; ES+: 736.54.

Example 360

(*rac.*)-(1*R**, 5*S**)-3-Acetyl-7-{4-[2-(2,3,5-trimethylphenoxy)ethyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid (2-chlorobenzyl)cyclopropylamide formate salt

Synthesized according to typical procedures A and E from bicyclononene AL38 and acetyl chloride. LC-MS: R_t = 0.93; ES+: 612.52.

30

Example 361

(*rac.*)-(1*R*^{*, 5*S*^{*})-3-Acetyl-7-{4-[2-(2,3,5-trimethylphenoxy)ethyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid (2-chlorobenzyl)cethylamide formate salt}

5 Synthesized according to typical procedures A and E from bicyclononene AL40 and acetyl chloride. LC-MS: R_t = 0.93; ES+: 600.50.

Example 362

10 **(*rac.*)-(1*R*^{*, 5*S*^{*})-3-Acetyl-7-{4-[2-(2,3,5-trimethylphenoxy)ethyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid (2-fluorobenzyl)cyclopropylamide formate salt}**

15 Synthesized according to typical procedures A and E from bicyclononene AL41 and acetyl chloride. LC-MS: R_t = 0.91; ES+: 596.53.

Example 363

20 **(*rac.*)-(1*R*^{*, 5*S*^{*})-3-Acetyl-7-{4-[2-(2,3,5-trimethylphenoxy)ethyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid (3-trifluoromethylbenzyl)cyclopropylamide formate salt}**

25 Synthesized according to typical procedures A and E from bicyclononene AL42 and acetyl chloride. LC-MS: R_t = 0.95; ES+: 646.54.

30

Example 364

(*rac.*)-(1*R*^{*, 5*S*^{*})-3-Acetyl-7-{4-[2-(2,3,5-trimethylphenoxy)ethyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid (2-methylbenzyl)cyclopropylamide formate salt}

Synthesized according to typical procedures A and E from bicyclononene AL43 and acetyl chloride. LC-MS: R_t = 0.93; ES+: 592.56.

Example 365

5

(rac.)-(1R, 5S*)-3-Acetyl-7-{4-[2-(2,3,5-trimethylphenoxy)ethyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid cyclopropyl-[2-(4-methoxyphenoxy)ethyl]amide formate salt*

10 Synthesized according to typical procedures A and E from bicyclononene AL44 and acetyl chloride. LC-MS: R_t = 0.92; ES+: 638.58.

Example 366

15 *(rac.)-5-((1R*, 5S*)-6-[(2-Chlorobenzyl)cyclopropylcarbamoyl]-7-{4-[2-(2,3,5-trimethylphenoxy)ethyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-en-3-yl)-5-oxo-pentanoic acid formate salt*

20 Synthesized according to typical procedures K and E from bicyclononene AL38 and glutaric anhydride. LC-MS: R_t = 0.91; ES+: 684.54.

Example 367

25 *(rac.)-5-((1R*, 5S*)-6-(Benzylcyclopropylcarbamoyl)-7-{4-[2-(2,3,5-trimethylphenoxy)ethyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-en-3-yl)-5-oxopentanoic acid formate salt*

Synthesized according to typical procedures K and E from bicyclononene AL39 and glutaric anhydride. LC-MS: R_t = 0.89; ES+: 650.57.

30

Example 368

(rac.)-5-((1R, 5S*)-6-[(2-Chlorobenzyl)ethylcarbamoyl]-7-{4-[2-(2,3,5-trimethylphenoxy)ethyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-en-3-yl)-5-oxopentanoic acid formate salt*

- 5 Synthesized according to typical procedures K and E from bicyclononene **AL40** and glutaric anhydride. LC-MS: R_t = 0.90; ES+: 672.55.

Example 369

- 10 *(rac.)-5-((1R*, 5S*)-6-[(2-Fluorobenzyl)cyclopropylcarbamoyl]-7-{4-[2-(2,3,5-trimethylphenoxy)ethyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-en-3-yl)-5-oxo-pentanoic acid formate salt*

15 Synthesized according to typical procedures K and E from bicyclononene **AL41** and glutaric anhydride. LC-MS: R_t = 0.89; ES+: 668.57.

Example 370

- 20 *(rac.)-5-((1R*, 5S*)-6-[(3-Trifluoromethylbenzyl)cyclopropylcarbamoyl]-7-{4-[2-(2,3,5-trimethylphenoxy)ethyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-en-3-yl)-5-oxo-pentanoic acid formate salt*

25 Synthesized according to typical procedures K and E from bicyclononene **AL42** and glutaric anhydride. LC-MS: R_t = 0.92; ES+: 718.52.

Example 371

- 30 *(rac.)-5-((1R*, 5S*)-6-[(2-Methylbenzyl)cyclopropylcarbamoyl]-7-{4-[2-(2,3,5-trimethylphenoxy)ethyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-en-3-yl)-5-oxo-pentanoic acid formate salt*

Synthesized according to typical procedures K and E from bicyclononene **AL43** and glutaric anhydride. LC-MS: R_t = 0.90; ES+: 664.59.

Example 372

5

(*rac.*)-5-((*1R*^{*,} *5S*^{*)}-6-{Cyclopropyl-[2-(4-methoxyphenoxy)ethyl]carbamoyl}-7-{4-[2-(2,3,5-trimethylphenoxy)ethyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-en-3-yl)-5-oxopentanoic acid formate salt

10 Synthesized according to typical procedures K and E from bicyclononene **AL44** and glutaric anhydride. LC-MS: R_t = 0.89; ES+: 710.56.

Example 373

15 (*rac.*)-5-((*1R*^{*,} *5S*^{*)}-6-{Cyclopropyl-[2-(3-methoxyphenoxy)ethyl]carbamoyl}-7-{4-[2-(2,3,5-trimethylphenoxy)ethyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-en-3-yl)-5-oxopentanoic acid formate salt

20 Synthesized according to typical procedures K and E from bicyclononene **AL45** and glutaric anhydride. LC-MS: R_t = 0.90; ES+: 710.54.

Example 374

25 (*rac.*)-5-((*1R*^{*,} *5S*^{*)}-6-{Cyclopropyl-[2-(3-methylphenoxy)ethyl]carbamoyl}-7-{4-[2-(2,3,5-trimethylphenoxy)ethyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-en-3-yl)-5-oxopentanoic acid formate salt

Synthesized according to typical procedures K and E from bicyclononene **AL46** and glutaric anhydride. LC-MS: R_t = 0.92; ES+: 694.57.

30

Example 375

(rac.)-5-((1R, 5S*)-6-[(2-Chlorobenzyl)ethylcarbamoyl]-7-{4-[2-(2,3,5-trimethylphenoxy)ethyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-en-3-yl)-5-oxopentanoic acid methyl ester formate salt*

5 Synthesized according to typical procedures A and E from bicyclononene AL40 and glutaric acid monomethylester chloride. LC-MS: R_t = 0.96; ES+: 686.54.

Example 376

10 *(rac.)-5-((1R*, 5S*)-6-[Cyclopropyl-(2-fluorobenzyl)carbamoyl]-7-{4-[2-(2,3,5-trimethylphenoxy)ethyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-en-3-yl)-5-oxopentanoic acid methyl ester*

15 Synthesized according to typical procedures A and E from bicyclononene AL41 and glutaric acid monomethylester chloride. LC-MS: R_t = 0.94; ES+: 682.57.

Example 377

20 *(rac.)-5-((1R*, 5S*)-6-[Cyclopropyl-(2-methylbenzyl)carbamoyl]-7-{4-[2-(2,3,5-trimethylphenoxy)ethyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-en-3-yl)-5-oxopentanoic acid methyl ester formate salt*

Synthesized according to typical procedures A and E from bicyclononene AL43 and glutaric acid monomethylester chloride. LC-MS: R_t = 0.96; ES+: 678.61.

25

Example 378

30 *(rac.)-5-((1R*, 5S*)-6-[(2-Chlorobenzyl)cyclopropylcarbamoyl]-7-{4-[2-(2,3,5-trimethylphenoxy)ethyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-en-3-yl)-2,2-dimethyl-5-oxopentanoic acid formate salt*

Synthesized according to typical procedures K and E from bicyclononene AL38 and 2,2-dimethylglutaric anhydride. LC-MS: R_t = 0.94; ES+: 712.51.

Example 379

5

(*rac.*)-5-((*1R*^{*}, *5S*^{*})-6-(Benzylcyclopropylcarbamoyl)-7-{4-[2-(2,3,5-trimethylphenoxy)ethyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-en-3-yl)-2,2-dimethyl-5-oxopentanoic acid formate salt

10 Synthesized according to typical procedures K and E from bicyclononene AL39 and 2,2-dimethylglutaric anhydride. LC-MS: R_t = 0.93; ES+: 678.60.

Example 380

15 (*rac.*)-5-((*1R*^{*}, *5S*^{*})-6-[(2-Chlorobenzyl)ethylcarbamoyl]-7-{4-[2-(2,3,5-trimethylphenoxy)ethyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-en-3-yl)-2,2-dimethyl-5-oxopentanoic acid formate salt

20 Synthesized according to typical procedures K and E from bicyclononene AL40 and 2,2-dimethylglutaric anhydride. LC-MS: R_t = 0.94; ES+: 700.52.

Example 381

25 (*rac.*)-5-((*1R*^{*}, *5S*^{*})-6-[Cyclopropyl-(2-fluorobenzyl)carbamoyl]-7-{4-[2-(2,3,5-trimethylphenoxy)ethyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-en-3-yl)-2,2-dimethyl-5-oxopentanoic acid formate salt

Synthesized according to typical procedures K and E from bicyclononene AL41 and 2,2-dimethylglutaric anhydride. LC-MS: R_t = 0.93; ES+: 696.57.

30

Example 382

(*rac.*)-5-((*1R*^{*}, *5S*^{*})-6-[Cyclopropyl-(2-methylbenzyl)carbamoyl]-7-{4-[2-(2,3,5-trimethylphenoxy)ethyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-en-3-yl)-2,2-dimethyl-5-oxopentanoic acid formate salt

5 Synthesized according to typical procedures K and E from bicyclononene AL43 and 2,2-dimethylglutaric anhydride. LC-MS: R_t = 0.94; ES+: 692.60.

Example 383

10 1:1-Mixture of (*rac.*)-(3*R*^{*})-4-((*1R*^{*}, *5S*^{*})-6-[cyclopropyl-(2-methylbenzyl)-carbamoyl]-7-{4-[2-(2,3,5-trimethylphenoxy)ethyl]phenyl}-3,9-diazabicyclo-[3.3.1]non-6-en-3-yl)-2,3-dihydroxy-4-oxobutyric acid formate salt and (*rac.*)-(3*S*^{*})-4-((*1R*^{*}, *5S*^{*})-6-[cyclopropyl-(2-methylbenzyl)carbamoyl]-7-{4-[2-(2,3,5-trimethylphenoxy)ethyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-en-3-yl)-2,3-dihydroxy-4-oxobutyric acid formate salt

15

2,3-dihydroxy-4-oxobutyric acid formate salt

Synthesized according to typical procedures K and E from bicyclononene AL43 and (S,R)-2,3-dihydroxysuccinic anhydride. LC-MS: R_t = 0.87; ES+: 682.57.

20 Example 384

(*rac.*)-(*1R*^{*}, *5S*^{*})-3-(4-Carbamoylbutyryl)-7-{4-[2-(2,3,5-trimethylphenoxy)-ethyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid benzyl-cyclopropylamide formate salt

25

Synthesized according to typical procedures G and E from bicyclononene AL39 and 4-carbamoylbutyric acid. LC-MS: R_t = 0.87; ES+: 649.59.

Example 385

30

(rac.)-(1R, 5S*)-3-(4-Carbamoylbutyryl)-7-{4-[2-(2,3,5-trimethylphenoxy)-ethyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid (2-chlorobenzyl)ethylamide formate salt*

5 Synthesized according to typical procedures G and E from bicyclononene AL40 and 4-carbamoylbutyric acid. LC-MS: R_t = 0.88; ES+: 671.56.

Example 386

10 *(rac.)-(1R*, 5S*)-3-(4-Carbamoylbutyryl)-7-{4-[2-(2,3,5-trimethylphenoxy)-ethyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid cyclopropyl-(2-fluorobenzyl)amide formate salt*

15 Synthesized according to typical procedures G and E from bicyclononene AL41 and 4-carbamoylbutyric acid. LC-MS: R_t = 0.87; ES+: 667.6.

Example 387

20 *(rac.)-(1R*, 5S*)-7-{4-[3-(2-Bromo-5-fluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid (2-chlorobenzyl)ethylamide formate salt*

25 Synthesized according to typical procedure E from bicyclononene AK22. LC-MS: R_t = 0.84; ES+: 628.36.

Example 388

30 *7-{4-[3-(2-Bromo-5-fluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]-non-6-ene-6-carboxylic acid cyclopropyl-(3,5-dimethoxybenzyl)amide formate salt*

Synthesized according to typical procedure E from bicyclononene AK35. LC-MS: R_t = 0.84; ES+: 666.43.

Example 389

5

7-{4-[3-(2-Bromo-5-fluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]-non-6-ene-6-carboxylic acid cyclopropyl-(2-methylbenzyl)amide formate salt

Synthesized according to typical procedure E from bicyclononene AK25. LC-MS: R_t = 0.85; ES+: 618.41.

Example 390

15 **7-{4-[3-(2-Bromo-5-fluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]-non-6-ene-6-carboxylic acid cyclopropyl-[2-(4-methoxyphenoxy)ethyl]amide formate salt**

Synthesized according to typical procedure E from bicyclononene AK26. LC-MS: R_t = 0.84; ES+: 664.43.

20

Example 391

25 **1:1-Mixture of (1*R*, 5*S*)-7-{4-[3-(2-bromo-5-fluorophenoxy)propyl]phenyl}-3-((2*S*, 4*R*)-4-hydroxypyrrolidine-2-carbonyl)-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid cyclopropyl-(3,5-dimethoxybenzyl)amide formate salt and (1*S*, 5*R*)-7-{4-[3-(2-bromo-5-fluorophenoxy)propyl]phenyl}-3-((2*S*, 4*R*)-4-hydroxypyrrolidine-2-carbonyl)-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid cyclopropyl-(3,5-dimethoxybenzyl)amide formate salt**

30 Synthesized according to typical procedures G, E and L from bicyclononene AL35 and BOC-L-hydroxyproline. LC-MS: R_t = 0.80; ES+: 777.50.

Example 392

- (*rac.*)-5-<{(1*R*^{*}, 5*S*^{*})-7-{4-[3-(2-Bromo-5-fluorophenoxy)propyl]phenyl}-6-[(2-chlorobenzyl)cyclopropylcarbamoyl]-3,9-diazabicyclo[3.3.1]non-6-en-3-yl}-5-oxopentanoic acid formate salt

Synthesized according to typical procedures K and E from bicyclononene **AL20** and glutaric anhydride. LC-MS: R_t = 0.89; ES+: 752.43.

10 **Example 393**

- (*rac.*)-5-<{(1*R*^{*}, 5*S*^{*})-7-{4-[3-(2-Bromo-5-fluorophenoxy)propyl]phenyl}-6-[(2-chlorobenzyl)cyclopropylcarbamoyl]-3,9-diazabicyclo[3.3.1]non-6-en-3-yl}-5-oxopentanoic acid methyl ester formate salt

15 Synthesized according to typical procedures A and E from bicyclononene **AL20** and glutaric acid monomethylester chloride. LC-MS: R_t = 0.95; ES+: 766.42.

20 **Example 394**

- (*rac.*)-5-<{(1*R*^{*}, 5*S*^{*})-7-{4-[3-(2-Bromo-5-fluorophenoxy)propyl]phenyl}-6-[(2-chlorobenzyl)cyclopropylcarbamoyl]-3,9-diazabicyclo[3.3.1]non-6-en-3-yl}-2,2-dimethyl-5-oxopentanoic acid formate salt

25 Synthesized according to typical procedures K and E from bicyclononene **AL20** and 2,2-dimethylglutaric anhydride. LC-MS: R_t = 0.93; ES+: 780.46.

Example 395

- 30 1:1-Mixture of (*rac.*)-(3*R*^{*})-4-<{(1*R*^{*}, 5*S*^{*})-7-{4-[3-(2-bromo-5-fluorophenoxy)-propyl]phenyl}-6-[(2-chlorobenzyl)cyclopropylcarbamoyl]-3,9-diazabicyclo-[3.3.1]non-6-en-3-yl}-2,3-dihydroxy-4-oxobutyric acid formate salt and (*rac.*)-

(3*S**)-4-((*1R**, *5S**)-7-{4-[3-(2-bromo-5-fluorophenoxy)propyl]phenyl}-6-[(2-chlorobenzyl)cyclopropylcarbamoyl]-3,9-diazabicyclo[3.3.1]non-6-en-3-yl)-2,3-dihydroxy-4-oxobutyric acid formate salt

- 5 Synthesized according to typical procedures K and E from bicyclononene AL20 and *meso*-dihydroxysuccinic anhydride. LC-MS: R_t = 0.87; ES+: 770.37.

Example 396

- 10 (*rac.*)-5-((*1R**, *5S**)-4-[3-(2-Bromo-5-fluorophenoxy)propyl]phenyl)-3-(4-carbamoylbutyryl)-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid (2-chlorobenzyl)cyclopropylamide formate salt

Synthesized according to typical procedures G and E from bicyclononene AL20 and 4-carbamoylbutyric acid. LC-MS: R_t = 0.88; ES+: 751.44.

Example 397

- 20 (*rac.*)-5-((*1R**, *5S**)-6-(Benzylcyclopropylcarbamoyl)-7-{4-[3-(2-bromo-5-fluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-en-3-yl)-5-oxo-pentanoic acid formate salt

Synthesized according to typical procedures K and E from bicyclononene AL21 and glutaric anhydride. LC-MS: R_t = 0.88; ES+: 718.46.

25

Example 398

- 30 (*rac.*)-5-((*1R**, *5S**)-6-(Benzylcyclopropylcarbamoyl)-7-{4-[3-(2-bromo-5-fluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-en-3-yl)-5-oxo-pentanoic acid methyl ester formate salt

Synthesized according to typical procedures A and E from bicyclononene **AL21** and glutaric acid monomethylester chloride. LC-MS: R_t = 0.93; ES+: 732.49.

Example 399

5

1:1-Mixture of (*rac.*)-(2*R*^{*}, 3*S*^{*})-4-((1*R*^{*}, 5*S*^{*})-6-(benzylcyclopropylcarbamoyl)-7-{4-[3-(2-bromo-5-fluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-en-3-yl)-2,3-dihydroxy-4-oxobutyric acid formate salt and (*rac.*)-(2*S*^{*}, 3*R*^{*})-4-((1*R*^{*}, 5*S*^{*})-6-(benzylcyclopropylcarbamoyl)-7-{4-[3-(2-bromo-5-fluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-en-3-yl)-2,3-dihydroxy-4-oxobutyric acid formate salt

10

Synthesized according to typical procedures K and E from bicyclononene **AL21** and (S,R)-2,3-dihydroxysuccinic anhydride. LC-MS: R_t = 0.86; ES+: 736.43.

15

Example 400

(*rac.*)-(1*R*^{*}, 5*S*^{*})-7-{4-[3-(2-Bromo-5-fluorophenoxy)propyl]phenyl}-3-(4-carbamoylbutyryl)-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid benzylcyclopropylamide formate salt

20

Synthesized according to typical procedures G and E from bicyclononene **AL21** and 4-carbamoylbutyric acid. LC-MS: R_t = 0.86; ES+: 717.46.

25

Example 401

1:1-Mixture of (*rac.*)-5-{(1*R*^{*}, 5*S*^{*})-7-{4-[3-(2-bromo-5-fluorophenoxy)propyl]phenyl}-6-[(2*R*^{*})-2-hydroxy-2-phenylethyl)methylcarbamoyl]-3,9-diazabicyclo[3.3.1]non-6-en-3-yl}-5-oxopentanoic acid formate salt and (*rac.*)-5-{(1*R*^{*}, 5*S*^{*})-7-{4-[3-(2-bromo-5-fluorophenoxy)propyl]phenyl}-6-[(2*S*^{*})-2-hydroxy-2-phenylethyl)methylcarbamoyl]-3,9-diazabicyclo[3.3.1]non-6-en-3-yl}-5-oxopentanoic acid formate salt

Synthesized according to typical procedures K and E from bicyclononene AL34 and glutaric anhydride. LC-MS: R_t = 0.85 ; ES+: 722.46

5 **Example 402**

1:1-Mixture of (rac.)-5-{(1R*, 5S*)-7-{4-[3-(2-bromo-5-fluorophenoxy)-propyl]phenyl}-6-[((2R*)-2-hydroxy-2-phenylethyl)methylcarbamoyl]-3,9-diazabicyclo[3.3.1]non-6-en-3-yl}-5-oxopentanoic acid methyl ester formate salt and (rac.)-5-{(1R*, 5S*)-7-{4-[3-(2-bromo-5-fluorophenoxy)propyl]-phenyl}-6-[((2S*)-2-hydroxy-2-phenylethyl)methylcarbamoyl]-3,9-diazabicyclo[3.3.1]non-6-en-3-yl}-5-oxopentanoic acid methyl ester formate salt

15 Synthesized according to typical procedures A and E from bicyclononene AL34 and glutaric acid monomethylester chloride. LC-MS: R_t = 0.88 ; ES+: 736.46

Example 403

1:1-Mixture of (rac.)-5-{(1R*, 5S*)-7-{4-[3-(2-bromo-5-fluorophenoxy)-propyl]phenyl}-6-[((2R*)-2-hydroxy-2-phenylethyl)methylcarbamoyl]-3,9-diazabicyclo[3.3.1]non-6-en-3-yl}-2,2-dimethyl-5-oxopentanoic acid formate salt and (rac.)-5-{(1R*, 5S*)-7-{4-[3-(2-bromo-5-fluorophenoxy)propyl]-phenyl}-6-[((2S*)-2-hydroxy-2-phenylethyl)methylcarbamoyl]-3,9-diazabicyclo[3.3.1]non-6-en-3-yl}-2,2-dimethyl-5-oxopentanoic acid formate salt

25 Synthesized according to typical procedures K and E from bicyclononene AL34 and 2,2-dimethylglutaric anhydride. LC-MS: R_t = 0.88 ; ES+: 750.47

Example 404

30 1:1-Mixture of (rac.)-(1R*, 5S*)-7-{4-[3-(2-bromo-5-fluorophenoxy)propyl]-phenyl}-3-(4-carbamoylbutyryl)-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxy-

lic acid ((2*R**)-2-hydroxy-2-phenylethyl)methylamide formate salt and (*rac.*)-(1*R**, 5*S**)-7-{4-[3-(2-bromo-5-fluorophenoxy)propyl]-phenyl}-3-(4-carbamoylbutyryl)-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid ((2*R**)-2-hydroxy-2-phenylethyl)methylamide formate salt

5

Synthesized according to typical procedures G and E from bicyclononene **AL34** and 4-carbamoylbutyric acid. LC-MS: R_t = 0.82 ; ES+: 721.46

Example 405

10

1:1-Mixture of (*rac.*)-(3*R**)-5-{(1*R**, 5*S**)-7-{4-[3-(2-bromo-5-fluorophenoxy)propyl]phenyl}-6-[(2-chlorobenzyl)cyclopropylcarbamoyl]-3,9-diazabicyclo[3.3.1]non-6-en-3-yl}-3-hydroxy-5-oxopentanoic acid formate salt and (*rac.*)-(3*S**)-5-{(1*R**, 5*S**)-7-{4-[3-(2-bromo-5-fluorophenoxy)propyl]phenyl}-6-[(2-chlorobenzyl)cyclopropylcarbamoyl]-3,9-diazabicyclo[3.3.1]non-6-en-3-yl}-3-hydroxy-5-oxopentanoic acid formate salt

15

Synthesized according to typical procedures K, E and L from bicyclononene **AL20** and 3-(*tert*-butyldimethylsilyloxy)glutaric anhydride. LC-MS: R_t = 0.88 ;

20 ES+: 768.40

Example 406

1:1-Mixture of (*rac.*)-(3*R**)-5-{(1*R**, 5*S**)-6-(benzylcyclopropylcarbamoyl)-7-{4-[3-(2-bromo-5-fluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-en-3-yl}-3-hydroxy-5-oxopentanoic acid formate salt and (*rac.*)-(3*S**)-5-{(1*R**, 5*S**)-6-(benzylcyclopropylcarbamoyl)-7-{4-[3-(2-bromo-5-fluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-en-3-yl}-3-hydroxy-5-oxopentanoic acid formate salt

25

30

Synthesized according to typical procedures K, E and L from bicyclononene **AL21** and 3-(*tert*-butyldimethylsilyloxy)glutaric anhydride. LC-MS: R_t = 0.86 ; ES+: 734.47

5 **Example 407**

(*rac.*)-(1*R*^{*}, 5*S*^{*})-7-{4-[2-(2-Chloro-4,5-dimethylphenoxy)ethoxy]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid [2-(2-chlorophenyl)ethyl]-cyclopropylamide formate salt

10

Synthesized according to typical procedure E from bicyclononene **AK65**. LC-MS: R_t = 0.84 ; ES+: 620.45

Example 408

15

(*rac.*)-(1*R*^{*}, 5*S*^{*})-7-{4-[2-(2-Chloro-4,5-dimethylphenoxy)ethoxy]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid cyclopropyl-(3,5-dimethoxybenzyl)amide formate salt

20

Synthesized according to typical procedure E from bicyclononene **AK70**. LC-MS: R_t = 0.84 ; ES+: 7632.32

Example 409

25

(*rac.*)-(1*R*^{*}, 5*S*^{*})-7-{4-[2-(2-Chloro-4,5-dimethylphenoxy)ethoxy]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid cyclopropyl-(*p*-toluyl-ethyl)amide formate salt

30

Synthesized according to typical procedure E from bicyclononene **AK71**. LC-MS: R_t = 0.85 ; ES+: 600.29

Example 410

(*rac.*)-(1*R*^{*}, 5*S*^{*})-3-Acetyl-7-{4-[2-(4-bromophenoxy)ethoxy]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid (2-chlorobenzyl)cyclopropylamide formate salt

5

Synthesized according to typical procedures A and E from bicyclononene AL73 and acetyl chloride. LC-MS: R_t = 0.90 ; ES+: 664.33

Example 411

10

(*rac.*)-(1*R*^{*}, 5*S*^{*})-3-Acetyl-7-{4-[2-(4-bromophenoxy)ethoxy]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid cyclopropyl-(3,5-dimethoxybenzyl)amide formate salt

15

Synthesized according to typical procedures A and E from bicyclononene AL85 and acetyl chloride. LC-MS: R_t = 0.88 ; ES+: 690.36

Example 412

20

(*rac.*)-5-((1*R*^{*}, 5*S*^{*})-7-{4-[2-(4-Bromophenoxy)ethoxy]phenyl}-6-[(2-chlorobenzyl)cyclopropylcarbamoyl]-3,9-diazabicyclo[3.3.1]non-6-en-3-yl)-5-oxopentanoic acid formate salt

25

Synthesized according to typical procedures K and E from bicyclononene AL73 and glutaric anhydride. LC-MS: R_t = 0.87 ; ES+: 736.31

Example 413

30

(*rac.*)-5-((1*R*^{*}, 5*S*^{*})-6-(Benzylcyclopropylcarbamoyl)-7-{4-[2-(4-bromophenoxy)ethoxy]phenyl}-3,9-diazabicyclo[3.3.1]non-6-en-3-yl)-5-oxopentanoic acid formate salt

Synthesized according to typical procedures K and E from bicyclononene AL74 and glutaric anhydride. LC-MS: R_t = 0.85 ; ES+: 702.33

Example 414

5

(*rac.*)-5-{(IR*, 5S*)}-7-{4-[2-(4-Bromophenoxy)ethoxy]phenyl}-6-[cyclopropyl-(2-fluorobenzyl)carbamoyl]-3,9-diazabicyclo[3.3.1]non-6-en-3-yl}-5-oxo-pentanoic acid formate salt

10 Synthesized according to typical procedures K and E from bicyclononene AL76 and glutaric anhydride. LC-MS: R_t = 0.86 ; ES+: 720.33

Example 415

15 (*rac.*)-5-{(IR*, 5S*)}-7-{4-[2-(4-Bromophenoxy)ethoxy]phenyl}-6-[cyclopropyl-(3-trifluoromethylbenzyl)carbamoyl]-3,9-diazabicyclo[3.3.1]non-6-en-3-yl}-5-oxopentanoic acid formate salt

20 Synthesized according to typical procedures K and E from bicyclononene AL77 and glutaric anhydride. LC-MS: R_t = 0.89 ; ES+: 770.27

Example 416

25 (*rac.*)-5-{(IR*, 5S*)}-7-{4-[2-(4-Bromophenoxy)ethoxy]phenyl}-6-[cyclopropyl-(2-methylbenzyl)carbamoyl]-3,9-diazabicyclo[3.3.1]non-6-en-3-yl}-5-oxo-pentanoic acid formate salt

Synthesized according to typical procedures K and E from bicyclononene AL78 and glutaric anhydride. LC-MS: R_t = 0.87 ; ES+: 716.34

30

Example 417

(*rac.*)-5-{(1*R*^{*}, 5*S*^{*})-7-[4-[2-(4-Bromophenoxy)ethoxy]phenyl}-6-[cyclopropyl[2-(3,4-dimethylphenoxy)ethyl]carbamoyl]-3,9-diazabicyclo[3.3.1]non-6-en-3-yl}-5-oxo-pentanoic acid formate salt

5 Synthesized according to typical procedures K and E from bicyclononene **AL80** and glutaric anhydride. LC-MS: R_t = 0.90 ; ES+: 760.37

Example 418

10 **(*rac.*)-5-{(1*R*^{*}, 5*S*^{*})-7-[4-[2-(4-Bromophenoxy)ethoxy]phenyl}-6-[2-(2-chlorophenyl)ethyl]cyclopropylcarbamoyl]-3,9-diazabicyclo[3.3.1]non-6-en-3-yl}-5-oxopentanoic acid formate salt**

15 Synthesized according to typical procedures K and E from bicyclononene **AL81** and glutaric anhydride. LC-MS: R_t = 0.88 ; ES+: 750.31

Example 419

20 **(*rac.*)-5-{(1*R*^{*}, 5*S*^{*})-7-[4-[2-(4-Bromophenoxy)ethoxy]phenyl}-6-[cyclopropyl[2-(4-fluorophenyl)ethyl]carbamoyl]-3,9-diazabicyclo[3.3.1]non-6-en-3-yl}-5-oxo-pentanoic acid formate salt**

Synthesized according to typical procedures K and E from bicyclononene **AL83** and glutaric anhydride. LC-MS: R_t = 0.87 ; ES+: 734.34

25

Example 420

30 **(*rac.*)-5-{(1*R*^{*}, 5*S*^{*})-7-[4-[2-(4-Bromophenoxy)ethoxy]phenyl}-6-[cyclopropyl-[2-(2-methylphenyl)ethyl]carbamoyl]-3,9-diazabicyclo[3.3.1]non-6-en-3-yl}-5-oxopentanoic acid formate salt**

Synthesized according to typical procedures K and E from bicyclononene **AL84** and glutaric anhydride. LC-MS: R_t = 0.88 ; ES+: 730.37

Example 421

5

(*rac.*)-5-{(1*R*^{*}, 5*S*^{*})-7-[4-[2-(4-Bromophenoxy)ethoxy]phenyl]-6-[cyclopropyl-(3,5-dimethoxybenzyl)carbamoyl]-3,9-diazabicyclo[3.3.1]non-6-en-3-yl}-5-oxopentanoic acid formate salt

10 Synthesized according to typical procedures K and E from bicyclononene **AL85** and glutaric anhydride. LC-MS: R_t = 0.86 ; ES+: 762.31

Example 422

15 (*rac.*)-5-{(1*R*^{*}, 5*S*^{*})-7-[4-[2-(4-Bromophenoxy)ethoxy]phenyl]-6-[cyclopropyl-[2-(4-methylphenyl)ethyl]carbamoyl]-3,9-diazabicyclo[3.3.1]non-6-en-3-yl}-5-oxopentanoic acid formate salt

Synthesized according to typical procedures K and E from bicyclononene **AL86**
20 and glutaric anhydride. LC-MS: R_t = 0.88 ; ES+: 730.37

Example 423

25 (*rac.*)-5-{(1*R*^{*}, 5*S*^{*})-7-[4-[2-(4-Bromophenoxy)ethoxy]phenyl]-6-[cyclopropyl-(2-fluorobenzyl)carbamoyl]-3,9-diazabicyclo[3.3.1]non-6-en-3-yl}-2,2-dimethyl-5-oxopentanoic acid formate salt

Synthesized according to typical procedures K and E from bicyclononene **AL76** and 2,2-dimethylglutaric anhydride. LC-MS: R_t = 0.89 ; ES+: 748.35

30

Example 424

(rac.)-5-((1R, 5S*)-7-{4-[2-(4-Bromophenoxy)ethoxy]phenyl}-6-{[2-(2-chlorophenyl)ethyl]cyclopropylcarbamoyl}-3,9-diazabicyclo[3.3.1]non-6-en-3-yl)-2,2-dimethyl-5-oxopentanoic acid formate salt*

- 5 Synthesized according to typical procedures K and E from bicyclononene AL81 and 2,2-dimethylglutaric anhydride. LC-MS: R_t = 0.91 ; ES+: 778.33

Example 425

- 10 *(rac.)-5-((1R*, 5S*)-7-{4-[2-(4-Bromophenoxy)ethoxy]phenyl}-6-{cyclopropyl-[2-(2,3-difluorophenyl)ethyl]carbamoyl}-3,9-diazabicyclo[3.3.1]non-6-en-3-yl)-2,2-dimethyl-5-oxopentanoic acid formate salt*

- 15 Synthesized according to typical procedures K and E from bicyclononene AL82 and 2,2-dimethylglutaric anhydride. LC-MS: R_t = 0.91 ; ES+: 780.36

Example 426

- 20 *(rac.)-5-((1R*, 5S*)-7-{4-[2-(4-Bromophenoxy)ethoxy]phenyl}-6-{cyclopropyl-[2-(2-methylphenyl)ethyl]carbamoyl}-3,9-diazabicyclo[3.3.1]non-6-en-3-yl)-2,2-dimethyl-5-oxopentanoic acid formate salt*

- 25 Synthesized according to typical procedures K and E from bicyclononene AL84 and 2,2-dimethylglutaric anhydride. LC-MS: R_t = 0.92 ; ES+: 7758.39

Example 427

- 30 *(rac.)-5-{(1R*, 5S*)-7-{4-[2-(4-Bromophenoxy)ethoxy]phenyl}-6-[cyclopropyl-(3,5-dimethoxybenzyl)carbamoyl]-3,9-diazabicyclo[3.3.1]non-6-en-3-yl}-2,2-dimethyl-5-oxopentanoic acid formate salt*

Synthesized according to typical procedures K and E from bicyclononene AL85 and 2,2-dimethylglutaric anhydride. LC-MS: R_t = 0.89 ; ES+: 790.39

Example 428

5

(*rac.*)-5-((*1R*^{*}, *5S*^{*})-7-{4-[2-(4-Bromophenoxy)ethoxy]phenyl}-6-{cyclopropyl-[2-(4-methylphenyl)ethyl]carbamoyl}-3,9-diazabicyclo[3.3.1]non-6-en-3-yl)-2,2-dimethyl-5-oxopentanoic acid formate salt

10 Synthesized according to typical procedures K and E from bicyclononene AL86 and 2,2-dimethylglutaric anhydride. LC-MS: R_t = 0.91 ; ES+: 758.38

Example 429

15 (*rac.*)-(*1R*^{*}, *5S*^{*})-7-{4-[2-(4-Bromophenoxy)ethoxy]phenyl}-3-(4-carbamoylbutyryl)-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid benzyl-cyclopropylamide formate salt

20 Synthesized according to typical procedures G and E from bicyclononene AL74 and 4-carbamoylbutyric acid. LC-MS: R_t = 0.85 ; ES+: 735.34

Example 430

25 (*rac.*)-(*1R*^{*}, *5S*^{*})-7-{4-[2-(4-Bromophenoxy)ethoxy]phenyl}-3-(4-carbamoylbutyryl)-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid (2-chlorobenzyl)-ethylamide formate salt

Synthesized according to typical procedures G and E from bicyclononene AL75 and 4-carbamoylbutyric acid. LC-MS: R_t = 0.85 ; ES+: 723.32

30

Example 431

(*rac.*)-(1*R*^{*,} 5*S*^{*})-7-{4-[2-(4-Bromophenoxy)ethoxy]phenyl}-3-(4-carbamoylbutyryl)-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid cyclopropyl-(2-fluorobenzyl)amide formate salt

5 Synthesized according to typical procedures G and E from bicyclononene AL76 and 4-carbamoylbutyric acid. LC-MS: R_t = 0.84 ; ES+: 719.33

Example 432

10 **(*rac.*)-(1*R*^{*,} 5*S*^{*})-7-{4-[2-(4-Bromophenoxy)ethoxy]phenyl}-3-(4-carbamoylbutyryl)-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid cyclopropyl-[2-(4-methoxyphenoxy)ethyl]amide formate salt**

15 Synthesized according to typical procedures G and E from bicyclononene AL79 and 4-carbamoylbutyric acid. LC-MS: R_t = 0.84 ; ES+: 761.34

Example 433

20 **(*rac.*)-(1*R*^{*,} 5*S*^{*})-7-{4-[2-(4-Bromophenoxy)ethoxy]phenyl}-3-(4-carbamoylbutyryl)-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid [2-(2-chlorophenyl)ethyl]cyclopropylamide formate salt**

25 Synthesized according to typical procedures G and E from bicyclononene AL81 and 4-carbamoylbutyric acid. LC-MS: R_t = 0.86 ; ES+: 749.32

25

Example 434

30 **(*rac.*)-(1*R*^{*,} 5*S*^{*})-7-{4-[2-(4-Bromophenoxy)ethoxy]phenyl}-3-(4-carbamoylbutyryl)-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid cyclopropyl-(3,5-dimethoxybenzyl)amide formate salt**

Synthesized according to typical procedures G and E from bicyclononene **AL85** and 4-carbamoylbutyric acid. LC-MS: R_t = 0.84 ; ES+: 761.35

Example 435

5

(*rac.*)-(1*R*^{*,} 5*S*^{*})-7-{4-[2-(4-Bromophenoxy)ethoxy]phenyl}-3-(4-carbamoylbutyryl)-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid cyclopropyl-[2-(2-hydroxyethyl)benzyl]amide formate salt

10 Synthesized according to typical procedures G and E from bicyclononene **AL87** and 4-carbamoylbutyric acid. LC-MS: R_t = 0.80 ; ES+: 745.40

Example 436

15 (*rac.*)-(1*R*^{*,} 5*S*^{*})-7-{4-[2-(2-Chloro-4,5-dimethylphenoxy)ethoxy]phenyl}-3-formyl-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid (2-chlorobenzyl)-ethylamide formate salt

Synthesized according to typical procedure E from bicyclononene **AK57**.
20 Obtained as side-product after purification by HPLC. LC-MS: R_t = 0.91 ; ES+: 622.25

Example 437

25 (*rac.*)-(1*R*^{*,} 5*S*^{*})-7-{4-[2-(2-Chloro-4,5-dimethylphenoxy)ethoxy]phenyl}-3-formyl-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid cyclopropyl-(2-*o*-tolylethyl)amide formate salt

Synthesized according to typical procedure E from bicyclononene **AK68**.
30 Obtained as side-product after purification by HPLC. LC-MS: R_t = 0.92 ; ES+: 628.31

Example 438

(*rac.*)-(1*R*^{*}, 5*S*^{*})-7-{4-[2-(2-Chloro-4,5-dimethylphenoxy)ethoxy]phenyl}-3-formyl-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid cyclopropyl-(3,5-dimethoxybenzyl)amide formate salt

Synthesized according to typical procedure E from bicyclononene AK70. Obtained as side-product after purification by HPLC. LC-MS: R_t = 0.90 ; ES+: 660.30

10

Example 439

(*rac.*)-(1*R*^{*}, 5*S*^{*})-3-Acetyl-7-{4-[2-(2-chloro-4,5-dimethylphenoxy)ethoxy]-phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid (2-chlorobenzyl)-cyclopropylamide formate salt

Synthesized according to typical procedures A and E from bicyclononene AL56 and acetyl chloride. LC-MS: R_t = 0.93 ; ES+: 648.37

20 **Example 440**

(*rac.*)-(1*R*^{*}, 5*S*^{*})-3-Acetyl-7-{4-[2-(2-Chloro-4,5-dimethylphenoxy)ethoxy]-phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid (2-chlorobenzyl)-ethylamide formate salt

25

Synthesized according to typical procedures A and E from bicyclononene AL57 and acetyl chloride. LC-MS: R_t = 0.92 ; ES+: 636.37

Example 441

30

(*rac.*)-(1*R*^{*}, 5*S*^{*})-3-Acetyl-7-{4-[2-(2-chloro-4,5-dimethylphenoxy)ethoxy]-phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid cyclopropyl-(2-fluorobenzyl)amide formate salt

- 5 Synthesized according to typical procedures A and E from bicyclononene AL58 and acetyl chloride. LC-MS: R_t = 0.90 ; ES+: 632.41

Example 442

- 10 **(*rac.*)-(1*R*^{*}, 5*S*^{*})-3-Acetyl-7-{4-[2-(2-chloro-4,5-dimethylphenoxy)ethoxy]-phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid cyclopropyl-(2-methylbenzyl)amide formate salt**

15 Synthesized according to typical procedures A and E from bicyclononene AL60 and acetyl chloride. LC-MS: R_t = 0.91 ; ES+: 628.43

Example 443

- 20 **(*rac.*)-(1*R*^{*}, 5*S*^{*})-3-Acetyl-7-{4-[2-(2-chloro-4,5-dimethylphenoxy)ethoxy]-phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid cyclopropyl-(3,5-dimethoxybenzyl)amide formate salt**

Synthesized according to typical procedures A and E from bicyclononene AL70 and acetyl chloride. LC-MS: R_t = 0.91 ; ES+: 674.44

25

Example 444

- 30 **(*rac.*)-(1*R*^{*}, 5*S*^{*})-3-Acetyl-7-{4-[2-(2-chloro-4,5-dimethylphenoxy)ethoxy]-phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid cyclopropyl-(2-p-tolyethyl)amide formate salt**

Synthesized according to typical procedures A and E from bicyclononene **AL71** and acetyl chloride. LC-MS: R_t = 0.92 ; ES+: 642.44

Example 445

5

(*rac.*)-5-((1*R*^{*}, 5*S*^{*})-6-[(2-Chlorobenzyl)cyclopropylcarbamoyl]-7-{4-[2-(2-chloro-4,5-dimethylphenoxy)ethoxy]phenyl}-3,9-diazabicyclo[3.3.1]non-6-en-3-yl)-5-oxopentanoic acid formate salt

10 Synthesized according to typical procedures K and E from bicyclononene **AL56** and glutaric anhydride. LC-MS: R_t = 0.90 ; ES+: 720.36

Example 446

15 (*rac.*)-5-((1*R*^{*}, 5*S*^{*})-6-[(2-Chlorobenzyl)ethylcarbamoyl]-7-{4-[2-(2-chloro-4,5-dimethylphenoxy)ethoxy]phenyl}-3,9-diazabicyclo[3.3.1]non-6-en-3-yl)-5-oxo-pentanoic acid formate salt

20 Synthesized according to typical procedures K and E from bicyclononene **AL57** and glutaric anhydride. LC-MS: R_t = 0.89 ; ES+: 708.35

Example 447

25 (*rac.*)-5-((1*R*^{*}, 5*S*^{*})-7-{4-[2-(2-Chloro-4,5-dimethylphenoxy)ethoxy]phenyl}-6-[cyclopropyl-(2-fluorobenzyl)carbamoyl]-3,9-diazabicyclo[3.3.1]non-6-en-3-yl)-5-oxopentanoic acid formate salt

Synthesized according to typical procedures K and E from bicyclononene **AL58** and glutaric anhydride. LC-MS: R_t = 0.88 ; ES+: 704.37

30

Example 448

(*rac.*)-5-((*1R*^{*}, *5S*^{*})-7-{4-[2-(2-Chloro-4,5-dimethylphenoxy)ethoxy]phenyl}-6-[cyclopropyl-(3-trifluoromethylbenzyl)carbamoyl]-3,9-diazabicyclo[3.3.1]-non-6-en-3-yl}-5-oxopentanoic acid formate salt

5 Synthesized according to typical procedures K and E from bicyclononene AL59
and glutaric anhydride. LC-MS: R_t = 0.91 ; ES+: 745.38

Example 449

10 **(*rac.*)-5-((*1R*^{*}, *5S*^{*})-7-{4-[2-(2-Chloro-4,5-dimethylphenoxy)ethoxy]phenyl}-6-[cyclopropyl-(2-methylbenzyl)carbamoyl]-3,9-diazabicyclo[3.3.1]non-6-en-3-yl}-5-oxopentanoic acid formate salt**

15 Synthesized according to typical procedures K and E from bicyclononene AL60
and glutaric anhydride. LC-MS: R_t = 0.89 ; ES+: 700.41

Example 450

20 **(*rac.*)-5-((*1R*^{*}, *5S*^{*})-7-{4-[2-(2-Chloro-4,5-dimethylphenoxy)ethoxy]phenyl}-6-[cyclopropyl-[2-(4-methoxyphenoxy)ethyl]carbamoyl]-3,9-diazabicyclo-[3.3.1]non-6-en-3-yl}-5-oxopentanoic acid formate salt**

25 Synthesized according to typical procedures K and E from bicyclononene AL61
and glutaric anhydride. LC-MS: R_t = 0.88 ; ES+: 746.42

Example 451

30 **(*rac.*)-5-((*1R*^{*}, *5S*^{*})-7-{4-[2-(2-Chloro-4,5-dimethylphenoxy)ethoxy]phenyl}-6-[cyclopropyl-[2-(3-methoxyphenoxy)ethyl]carbamoyl]-3,9-diazabicyclo-[3.3.1]non-6-en-3-yl}-5-oxopentanoic acid formate salt**

Synthesized according to typical procedures K and E from bicyclononene AL62 and glutaric anhydride. LC-MS: R_t = 0.88 ; ES+: 746.41

Example 452

5

(*rac.*)-5-((1*R**, 5*S**)-7-{4-[2-(2-Chloro-4,5-dimethylphenoxy)ethoxy]phenyl}-6-{cyclopropyl-[2-(3-methylphenoxy)ethyl]carbamoyl}-3,9-diazabicyclo[3.3.1]-non-6-en-3-yl)-5-oxopentanoic acid formate salt

10 Synthesized according to typical procedures K and E from bicyclononene AL63 and glutaric anhydride. LC-MS: R_t = 0.91 ; ES+: 730.44

Example 453

15 (*rac.*)-5-[(1*R**, 5*S**)-7-{4-[2-(2-Chloro-4,5-dimethylphenoxy)ethoxy]phenyl}-6-(cyclopropylphenethylcarbamoyl)-3,9-diazabicyclo[3.3.1]non-6-en-3-yl]-5-oxo-pentanoic acid formate salt

20 Synthesized according to typical procedures K and E from bicyclononene AL64 and glutaric anhydride. LC-MS: R_t = 0.88 ; ES+: 700.41

Example 454

25 (*rac.*)-5-((1*R**, 5*S**)-7-{4-[2-(2-Chloro-4,5-dimethylphenoxy)ethoxy]phenyl}-6-{{[2-(2-chlorophenyl)ethyl]cyclopropyl-carbamoyl}-3,9-diazabicyclo[3.3.1]-non-6-en-3-yl)-5-oxopentanoic acid formate salt

Synthesized according to typical procedures K and E from bicyclononene AL65 and glutaric anhydride. LC-MS: R_t = 0.90 ; ES+: 734.39

30

Example 455

(rac.)-5-((1R, 5S*)-7-{4-[2-(2-Chloro-4,5-dimethylphenoxy)ethoxy]phenyl}-6-{cyclopropyl-[2-(2,3-difluorophenyl)ethyl]carbamoyl}-3,9-diazabicyclo[3.3.1]-non-6-en-3-yl)-5-oxopentanoic acid formate salt*

- 5 Synthesized according to typical procedures K and E from bicyclononene AL66 and glutaric anhydride. LC-MS: R_t = 0.90 ; ES+: 736.41

Example 456

- 10 *(rac.)-5-((1R*, 5S*)-7-{4-[2-(2-Chloro-4,5-dimethylphenoxy)ethoxy]phenyl}-6-{cyclopropyl-[2-(4-fluorophenyl)ethyl]carbamoyl}-3,9-diazabicyclo[3.3.1]-non-6-en-3-yl)-5-oxopentanoic acid formate salt*

15 Synthesized according to typical procedures K and E from bicyclononene AL67 and glutaric anhydride. LC-MS: R_t = 0.89 ; ES+: 718.40

Example 457

- 20 *(rac.)-5-((1R*, 5S*)-7-{4-[2-(2-Chloro-4,5-dimethylphenoxy)ethoxy]phenyl}-6-{cyclopropyl-[2-(2-methylphenyl)ethyl]carbamoyl}-3,9-diazabicyclo[3.3.1]-non-6-en-3-yl)-5-oxopentanoic acid formate salt*

25 Synthesized according to typical procedures K and E from bicyclononene AL68 and glutaric anhydride. LC-MS: R_t = 0.90 ; ES+: 714.43

Example 458

- 30 *(rac.)-5-((1R*, 5S*)-7-{4-[2-(2-Chloro-4,5-dimethylphenoxy)ethoxy]phenyl}-6-[cyclopropyl-(3,5-dimethoxybenzyl)carbamoyl]-3,9-diazabicyclo[3.3.1]non-6-en-3-yl)-5-oxopentanoic acid formate salt*

Synthesized according to typical procedures K and E from bicyclononene **AL70** and glutaric anhydride. LC-MS: R_t = 0.89 ; ES+: 746.41

Example 459

5

(*rac.*)-5-((*1R*^{*}, *5S*^{*})-7-{4-[2-(2-Chloro-4,5-dimethylphenoxy)ethoxy]phenyl}-6-{cyclopropyl-[2-(4-methylphenyl)ethyl]carbamoyl}-3,9-diazabicyclo[3.3.1]-non-6-en-3-yl)-5-oxopentanoic acid formate salt

10 Synthesized according to typical procedures K and E from bicyclononene **AL71** and glutaric anhydride. LC-MS: R_t = 0.90 ; ES+: 714.41

Example 460

15 (*rac.*)-5-((*1R*^{*}, *5S*^{*})-6-[(2-Chlorobenzyl)cyclopropylcarbamoyl]-7-{4-[2-(2-chloro-4,5-dimethylphenoxy)ethoxy]phenyl}-3,9-diazabicyclo[3.3.1]non-6-en-3-yl)-5-oxopentanoic acid methyl ester formate salt

20 Synthesized according to typical procedures A and E from bicyclononene **AL56** and glutaric acid monomethylester chloride. LC-MS: R_t = 0.95 ; ES+: 734.37

Example 461

25 (*rac.*)-5-((*1R*^{*}, *5S*^{*})-6-[(2-Chlorobenzyl)ethylcarbamoyl]-7-{4-[2-(2-chloro-4,5-dimethylphenoxy)ethoxy]phenyl}-3,9-diazabicyclo[3.3.1]non-6-en-3-yl)-5-oxopentanoic acid methyl ester formate salt

Synthesized according to typical procedures A and E from bicyclononene **AL57** and glutaric acid monomethylester chloride. LC-MS: R_t = 0.94 ; ES+: 722.37

30

Example 462

(rac.)-5-{(1R, 5S*)-7-{4-[2-(2-Chloro-4,5-dimethylphenoxy)ethoxy]phenyl}-6-[cyclopropyl-(2-fluorobenzyl)carbamoyl]-3,9-diazabicyclo[3.3.1]non-6-en-3-yl}-5-oxopentanoic acid methyl ester formate salt*

- 5 Synthesized according to typical procedures A and E from bicyclononene AL58 and glutaric acid monomethylester chloride. LC-MS: R_t = 0.93 ; ES+: 718.40

Example 463

- 10 *(rac.)-5-{(1R*, 5S*)-7-{4-[2-(2-Chloro-4,5-dimethylphenoxy)ethoxy]phenyl}-6-[cyclopropyl-(3-trifluoromethylbenzyl)carbamoyl]-3,9-diazabicyclo[3.3.1]-non-6-en-3-yl}-5-oxopentanoic acid methyl ester formate salt*

15 Synthesized according to typical procedures A and E from bicyclononene AL59 and glutaric acid monomethylester chloride. LC-MS: R_t = 0.96 ; ES+: 768.38

Example 464

- 20 *(rac.)-5-{(1R*, 5S*)-7-{4-[2-(2-Chloro-4,5-dimethylphenoxy)ethoxy]phenyl}-6-[cyclopropyl-(2-methylbenzyl)carbamoyl]-3,9-diazabicyclo[3.3.1]non-6-en-3-yl}-5-oxopentanoic acid methyl ester formate salt*

25 Synthesized according to typical procedures A and E from bicyclononene AL60 and glutaric acid monomethylester chloride. LC-MS: R_t = 0.94 ; ES+: 714.43

Example 465

- 30 *(rac.)-5-[(1R*, 5S*)-7-{4-[2-(2-Chloro-4,5-dimethylphenoxy)ethoxy]phenyl}-6-(cyclopropylphenethylcarbamoyl)-3,9-diazabicyclo[3.3.1]non-6-en-3-yl]-5-oxopentanoic acid methyl ester formate salt*

Synthesized according to typical procedures A and E from bicyclononene **AL64** and glutaric acid monomethylester chloride. LC-MS: R_t = 0.93 ; ES+: 714.41

Example 466

5

(*rac.*)-5-((*1R*^{*}, *5S*^{*})-7-{4-[2-(2-Chloro-4,5-dimethylphenoxy)ethoxy]phenyl}-6-{|[2-(2-chlorophenyl)ethyl]cyclopropylcarbamoyl}-3,9-diazabicyclo[3.3.1]non-6-en-3-yl)-5-oxopentanoic acid methyl ester formate salt

10 Synthesized according to typical procedures A and E from bicyclononene **AL65** and glutaric acid monomethylester chloride. LC-MS: R_t = 0.93 ; ES+: 748.39

Example 467

15 (*rac.*)-5-((*1R*^{*}, *5S*^{*})-7-{4-[2-(2-Chloro-4,5-dimethylphenoxy)ethoxy]phenyl}-6-{|[2-(2,3-difluorophenyl)ethyl]cyclopropylcarbamoyl}-3,9-diazabicyclo[3.3.1]-non-6-en-3-yl)-5-oxopentanoic acid methyl ester formate salt

20 Synthesized according to typical procedures A and E from bicyclononene **AL66** and glutaric acid monomethylester chloride. LC-MS: R_t = 0.94 ; ES+: 750.40

Example 468

25 (*rac.*)-5-((*1R*^{*}, *5S*^{*})-7-{4-[2-(2-Chloro-4,5-dimethylphenoxy)ethoxy]phenyl}-6-{|[2-(4-fluorophenyl)ethyl]cyclopropylcarbamoyl}-3,9-diazabicyclo[3.3.1]non-6-en-3-yl)-5-oxopentanoic acid methyl ester formate salt

Synthesized according to typical procedures A and E from bicyclononene **AL67** and glutaric acid monomethylester chloride. LC-MS: R_t = 0.93 ; ES+: 732.44

30

Example 469

(rac.)-5-((1R, 5S*)-7-{4-[2-(2-Chloro-4,5-dimethylphenoxy)ethoxy]phenyl}-6-[[2-(2-methylphenyl)ethyl]cyclopropylcarbamoyl]-3,9-diazabicyclo[3.3.1]non-6-en-3-yl)-5-oxopentanoic acid methyl ester formate salt*

- 5 Synthesized according to typical procedures A and E from bicyclononene **AL68** and glutaric acid monomethylester chloride. LC-MS: R_t = 0.95 ; ES+: 728.45

Example 470

- 10 *(rac.)-5-((1R*, 5S*)-7-{4-[2-(2-Chloro-4,5-dimethylphenoxy)ethoxy]phenyl}-6-[[cyclopropyl-(3,5-dimethoxybenzyl)carbamoyl]-3,9-diazabicyclo[3.3.1]non-6-en-3-yl)-5-oxopentanoic acid methyl ester formate salt*

Synthesized according to typical procedures A and E from bicyclononene **AL70**
15 and glutaric acid monomethylester chloride. LC-MS: R_t = 0.93 ; ES+: 760.40

Example 471

- 10 *(rac.)-5-((1R*, 5S*)-7-{4-[2-(2-Chloro-4,5-dimethylphenoxy)ethoxy]phenyl}-6-[[2-(4-methylphenyl)ethyl]cyclopropylcarbamoyl]-3,9-diazabicyclo[3.3.1]non-6-en-3-yl)-5-oxopentanoic acid methyl ester formate salt*

Synthesized according to typical procedures A and E from bicyclononene **AL71**
25 and glutaric acid monomethylester chloride. LC-MS: R_t = 0.95 ; ES+: 728.43

Example 472

- 10 *(rac.)-5-((1R*, 5S*)-6-[(2-Chlorobenzyl)cyclopropylcarbamoyl]-7-{4-[2-(2-chloro-4,5-dimethylphenoxy)ethoxy]phenyl}-3,9-diazabicyclo[3.3.1]non-6-en-3-yl)-2,2-dimethyl-5-oxopentanoic acid formate salt*

Synthesized according to typical procedures K and E from bicyclononene **AL56** and 2,2-dimethylglutaric anhydride. LC-MS: R_t = 0.93 ; ES+: 748.40

Example 473

5

(*rac.*)-5-((*1R*^{*}, *5S*^{*})-6-[(2-Chlorobenzyl)ethylcarbamoyl]-7-{4-[2-(2-chloro-4,5-dimethylphenoxy)ethoxy]phenyl}-3,9-diazabicyclo[3.3.1]non-6-en-3-yl)-2,2-dimethyl-5-oxopentanoic acid formate salt

10 Synthesized according to typical procedures K and E from bicyclononene **AL57** and 2,2-dimethylglutaric anhydride. LC-MS: R_t = 0.93 ; ES+: 736.42

Example 474

15 (*rac.*)-5-((*1R*^{*}, *5S*^{*})-7-{4-[2-(2-Chloro-4,5-dimethylphenoxy)ethoxy]phenyl}-6-[cyclopropyl-(2-fluorobenzyl)carbamoyl]-3,9-diazabicyclo[3.3.1]non-6-en-3-yl)-2,2-dimethyl-5-oxopentanoic acid formate salt

20 Synthesized according to typical procedures K and E from bicyclononene **AL58** and 2,2-dimethylglutaric anhydride. LC-MS: R_t = 0.91 ; ES+: 732.45

Example 475

25 (*rac.*)-5-((*1R*^{*}, *5S*^{*})-7-{4-[2-(2-Chloro-4,5-dimethylphenoxy)ethoxy]phenyl}-6-[cyclopropyl-(2-methylbenzyl)carbamoyl]-3,9-diazabicyclo[3.3.1]non-6-en-3-yl)-2,2-dimethyl-5-oxopentanoic acid formate salt

Synthesized according to typical procedures K and E from bicyclononene **AL60** and 2,2-dimethylglutaric anhydride. LC-MS: R_t = 0.93 ; ES+: 728.48

30

Example 476

(rac.)-5-[(1*R*^{*, 5*S*^{*})-7-{4-[2-(2-Chloro-4,5-dimethylphenoxy)ethoxy]phenyl}-6-(cyclopropylphenethylcarbamoyl)-3,9-diazabicyclo[3.3.1]non-6-en-3-yl]-2,2-dimethyl-5-oxopentanoic acid formate salt}

5 Synthesized according to typical procedures K and E from bicyclononene **AL64** and 2,2-dimethylglutaric anhydride. LC-MS: R_t = 0.92 ; ES+: 728.49

Example 477

10 **(rac.)-5-((1*R*^{*, 5*S*^{*})-7-{4-[2-(2-Chloro-4,5-dimethylphenoxy)ethoxy]phenyl}-6-{[2-(2-chlorophenyl)ethyl]cyclopropylcarbamoyl}-3,9-diazabicyclo[3.3.1]non-6-en-3-yl]-2,2-dimethyl-5-oxopentanoic acid formate salt}**

15 Synthesized according to typical procedures K and E from bicyclononene **AL65** and 2,2-dimethylglutaric anhydride. LC-MS: R_t = 0.93 ; ES+: 762.42

Example 478

20 **(rac.)-5-((1*R*^{*, 5*S*^{*})-7-{4-[2-(2-Chloro-4,5-dimethylphenoxy)ethoxy]phenyl}-6-{[2-(2,3-difluorophenyl)ethyl]cyclopropylcarbamoyl}-3,9-diazabicyclo[3.3.1]-non-6-en-3-yl]-2,2-dimethyl-5-oxopentanoic acid formate salt}**

25 Synthesized according to typical procedures K and E from bicyclononene **AL66** and 2,2-dimethylglutaric anhydride. LC-MS: R_t = 0.93 ; ES+: 764.44

Example 479

30 **(rac.)-5-((1*R*^{*, 5*S*^{*})-7-{4-[2-(2-Chloro-4,5-dimethylphenoxy)ethoxy]phenyl}-6-{[2-(4-fluorophenyl)ethyl]cyclopropylcarbamoyl}-3,9-diazabicyclo[3.3.1]non-6-en-3-yl]-2,2-dimethyl-5-oxopentanoic acid formate salt}**

Synthesized according to typical procedures K and E from bicyclononene AL67 and 2,2-dimethylglutaric anhydride. LC-MS: R_t = 0.92 ; ES+: 746.48

Example 480

5

(*rac.*)-5-((1*R*^{*, 5*S*^{*})-7-{4-[2-(2-Chloro-4,5-dimethylphenoxy)ethoxy]phenyl}-6- {[2-(2-methylphenyl)ethyl]cyclopropylcarbamoyl}-3,9-diazabicyclo[3.3.1]non-6-en-3-yl)-2,2-dimethyl-5-oxopentanoic acid formate salt}

10 Synthesized according to typical procedures K and E from bicyclononene AL68 and 2,2-dimethylglutaric anhydride. LC-MS: R_t = 0.93 ; ES+: 742.51

Example 481

15 (*rac.*)-5-((1*R*^{*, 5*S*^{*})-7-{4-[2-(2-Chloro-4,5-dimethylphenoxy)ethoxy]phenyl}-6- {[cyclopropyl-(3,5-dimethoxybenzyl)carbamoyl]-3,9-diazabicyclo[3.3.1]non-6-en-3-yl)-2,2-dimethyl-5-oxopentanoic acid formate salt}

20 Synthesized according to typical procedures K and E from bicyclononene AL70 and 2,2-dimethylglutaric anhydride. LC-MS: R_t = 0.92 ; ES+: 774.45

Example 482

25 (*rac.*)-5-((1*R*^{*, 5*S*^{*})-7-{4-[2-(2-Chloro-4,5-dimethylphenoxy)ethoxy]phenyl}-6- {[2-(4-methylphenyl)ethyl]cyclopropylcarbamoyl}-3,9-diazabicyclo[3.3.1]non-6-en-3-yl)-2,2-dimethyl-5-oxopentanoic acid formate salt}

Synthesized according to typical procedures K and E from bicyclononene AL71 and 2,2-dimethylglutaric anhydride. LC-MS: R_t = 0.93 ; ES+: 742.51

30

Example 483

(*rac.*)-(1*R*^{*}, 5*S*^{*})-3-(4-Carbamoylbutyryl)-7-{4-[2-(2-chloro-4,5-dimethyl-phenoxy)ethoxy]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid (2-chlorobenzyl)cyclopropylamide formate salt

5 Synthesized according to typical procedures G and E from bicyclononene AL56 and 4-carbamoylbutyric acid. LC-MS: R_t = 0.88 ; ES+: 719.43

Example 484

10 **(*rac.*)-(1*R*^{*}, 5*S*^{*})-3-(4-Carbamoylbutyryl)-7-{4-[2-(2-chloro-4,5-dimethyl-phenoxy)ethoxy]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid (2-chlorobenzyl)ethylamide formate salt**

15 Synthesized according to typical procedures G and E from bicyclononene AL57 and 4-carbamoylbutyric acid. LC-MS: R_t = 0.87 ; ES+: 707.39

Example 485

20 **(*rac.*)-(1*R*^{*}, 5*S*^{*})-3-(4-Carbamoylbutyryl)-7-{4-[2-(2-chloro-4,5-dimethyl-phenoxy)ethoxy]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid cyclopropyl-(2-fluorobenzyl)amide formate salt**

Synthesized according to typical procedures G and E from bicyclononene AL58 and 4-carbamoylbutyric acid. LC-MS: R_t = 0.86 ; ES+: 703.43

25

Example 486

30 **(*rac.*)-(1*R*^{*}, 5*S*^{*})-3-(4-Carbamoylbutyryl)-7-{4-[2-(2-chloro-4,5-dimethyl-phenoxy)ethoxy]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid cyclopropyl-(3-trifluoromethylbenzyl)amide formate salt**

Synthesized according to typical procedures G and E from bicyclononene **AL59** and 4-carbamoylbutyric acid. LC-MS: R_t = 0.89 ; ES+: 753.44

Example 487

5

(*rac.*)-(1*R*^{*}, 5*S*^{*})-3-(4-Carbamoylbutyryl)-7-{4-[2-(2-chloro-4,5-dimethyl-phenoxy)ethoxy]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid cyclopropyl-(2-methylbenzyl)amide formate salt

10 Synthesized according to typical procedures G and E from bicyclononene **AL60** and 4-carbamoylbutyric acid. LC-MS: R_t = 0.87 ; ES+: 699.47

Example 488

15 (*rac.*)-(1*R*^{*}, 5*S*^{*})-3-(4-Carbamoylbutyryl)-7-{4-[2-(2-chloro-4,5-dimethyl-phenoxy)ethoxy]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid cyclopropyl-[2-(4-methoxyphenoxy)ethyl]amide formate salt

20 Synthesized according to typical procedures G and E from bicyclononene **AL61** and 4-carbamoylbutyric acid. LC-MS: R_t = 0.86 ; ES+: 745.46

Example 489

25 (*rac.*)-(1*R*^{*}, 5*S*^{*})-3-(4-Carbamoylbutyryl)-7-{4-[2-(2-chloro-4,5-dimethyl-phenoxy)ethoxy]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid cyclopropylphenethylamide formate salt

30 Synthesized according to typical procedures G and E from bicyclononene **AL64** and 4-carbamoylbutyric acid. LC-MS: R_t = 0.86 ; ES+: 699.44

Example 490

(*rac.*)-(1*R*^{*}, 5*S*^{*})-3-(4-Carbamoylbutyryl)-7-{4-[2-(2-chloro-4,5-dimethyl-phenoxy)ethoxy]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid [2-(2-chlorophenyl)ethyl]cyclopropylamide formate salt

5 Synthesized according to typical procedures G and E from bicyclononene AL65 and 4-carbamoylbutyric acid. LC-MS: R_t = 0.88 ; ES+: 733.42

Example 491

10 **(*rac.*)-(1*R*^{*}, 5*S*^{*})-3-(4-Carbamoylbutyryl)-7-{4-[2-(2-chloro-4,5-dimethyl-phenoxy)ethoxy]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid cyclopropyl-[2-(2,3-difluorophenyl)ethyl]amide formate salt**

15 Synthesized according to typical procedures G and E from bicyclononene AL66 and 4-carbamoylbutyric acid. LC-MS: R_t = 0.88 ; ES+: 735.43

Example 492

20 **(*rac.*)-(1*R*^{*}, 5*S*^{*})-3-(4-Carbamoylbutyryl)-7-{4-[2-(2-chloro-4,5-dimethyl-phenoxy)ethoxy]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid cyclopropyl-[2-(4-fluorophenyl)ethyl]amide formate salt**

25 Synthesized according to typical procedures G and E from bicyclononene AL67 and 4-carbamoylbutyric acid. LC-MS: R_t = 0.87 ; ES+: 717.43

Example 493

30 **(*rac.*)-(1*R*^{*}, 5*S*^{*})-3-(4-Carbamoylbutyryl)-7-{4-[2-(2-chloro-4,5-dimethyl-phenoxy)ethoxy]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid cyclopropyl-(2-*o*-tolylethyl)amide formate salt**

Synthesized according to typical procedures G and E from bicyclononene AL68 and 4-carbamoylbutyric acid. LC-MS: R_t = 0.88 ; ES+: 713.46

Example 494

5

(rac.)-(1R, 5S*)-3-(4-Carbamoylbutyryl)-7-{4-[2-(2-chloro-4,5-dimethylphenoxy)ethoxy]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid cyclopropyl-(3,5-dimethoxybenzyl)amide formate salt*

10 Synthesized according to typical procedures G and E from bicyclononene AL70 and 4-carbamoylbutyric acid. LC-MS: R_t = 0.87 ; ES+: 745.45

Example 495

15 *(rac.)-(1R*, 5S*)-3-(4-Carbamoylbutyryl)-7-{4-[2-(2-chloro-4,5-dimethylphenoxy)ethoxy]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid cyclopropyl-(2-p-tolylethyl)amide formate salt*

20 Synthesized according to typical procedures G and E from bicyclononene AL71 and 4-carbamoylbutyric acid. LC-MS: R_t = 0.88 ; ES+: 713.46

Example 496

25 *1:1-Mixture of (rac.)-(3R*)-5-((1R*, 5S*)-6-[(2-chlorobenzyl)cyclopropylcarbamoyl]-7-{4-[2-(2-chloro-4,5-dimethylphenoxy)ethoxy]phenyl}-3,9-diazabicyclo[3.3.1]non-6-en-3-yl)-3-hydroxy-5-oxopentanoic acid formate salt and (rac.)-(3S*)-5-((1R*, 5S*)-6-[(2-chlorobenzyl)cyclopropyl-carbamoyl]-7-{4-[2-(2-chloro-4,5-dimethylphenoxy)ethoxy]phenyl}-3,9-diazabicyclo[3.3.1]non-6-en-3-yl)-3-hydroxy-5-oxopentanoic acid formate salt*

30

Synthesized according to typical procedures K, E and L from bicyclononene AL56 and 3-(*tert*-butyldimethylsilyloxy)glutaric anhydride. LC-MS: R_t = 0.88 ; ES+: 736.43

5 **Example 497**

1:1-Mixture of (*rac.*)-(3*R**)-5-((1*R**, 5*S**)-6-[(2-chlorobenzyl)ethylcarbamoyl]-7-{4-[2-(2-chloro-4,5-dimethylphenoxy)ethoxy]phenyl}-3,9-diazabicyclo[3.3.1]-non-6-en-3-yl)-3-hydroxy-5-oxopentanoic acid formate salt and
10 (*rac.*)-(3*S**)-5-((1*R**, 5*S**)-6-[(2-chlorobenzyl)ethylcarbamoyl]-7-{4-[2-(2-chloro-4,5-dimethylphenoxy)ethoxy]phenyl}-3,9-diazabicyclo[3.3.1]-non-6-en-3-yl)-3-hydroxy-5-oxopentanoic acid formate salt

Synthesized according to typical procedures K, E and L from bicyclononene
15 AL57 and 3-(*tert*-butyldimethylsilyloxy)glutaric anhydride. LC-MS: R_t = 0.87 ;
ES+: 724.40

Example 498

20 1:1-Mixture of (*rac.*)-(3*R**)-5-((1*R**, 5*S**)-7-{4-[2-(2-chloro-4,5-dimethylphenoxy)ethoxy]phenyl}-6-[cyclopropyl-(2-fluorobenzyl)carbamoyl]-3,9-diazabicyclo[3.3.1]non-6-en-3-yl)-3-hydroxy-5-oxopentanoic acid formate salt and (*rac.*)-(3*S**)-5-((1*R**, 5*S**)-7-{4-[2-(2-chloro-4,5-dimethylphenoxy)ethoxy]phenyl}-6-[cyclopropyl-(2-fluorobenzyl)carbamoyl]-3,9-diazabicyclo[3.3.1]non-6-en-3-yl)-3-hydroxy-5-oxopentanoic acid formate salt
25

Synthesized according to typical procedures K, E and L from bicyclononene AL58 and 3-(*tert*-butyldimethylsilyloxy)glutaric anhydride. LC-MS: R_t = 0.86 ;
ES+: 720.43

Example 499

1:1-Mixture of (*rac.*)-(3*R**)-5-{(*IR**, *5S**)-7-{4-[2-(2-chloro-4,5-dimethylphenoxy)ethoxy]phenyl}-6-[cyclopropyl-(3-trifluoromethylbenzyl)-carbamoyl]-3,9-diazabicyclo[3.3.1]non-6-en-3-yl}-3-hydroxy-5-oxopentanoic acid formate salt and (*rac.*)-(3*S**)-5-{(*IR**, *5S**)-7-{4-[2-(2-chloro-4,5-dimethylphenoxy)ethoxy]phenyl}-6-[cyclopropyl-(3-trifluoromethylbenzyl)-carbamoyl]-3,9-diazabicyclo[3.3.1]non-6-en-3-yl}-3-hydroxy-5-oxopentanoic acid formate salt

Synthesized according to typical procedures K, E and L from bicyclononene AL59 and 3-(*tert*-butyldimethylsilyloxy)glutaric anhydride. LC-MS: R_t = 0.89 ; ES+: 770.40

Example 500

1:1-Mixture of (*rac.*)-(3*R**)-5-{(*IR**, *5S**)-7-{4-[2-(2-chloro-4,5-dimethylphenoxy)ethoxy]phenyl}-6-[cyclopropyl-[2-(4-methoxyphenoxy)ethyl]-carbamoyl]-3,9-diazabicyclo[3.3.1]non-6-en-3-yl}-3-hydroxy-5-oxopentanoic acid formate salt and (*rac.*)-(3*S**)-5-{(*IR**, *5S**)-7-{4-[2-(2-chloro-4,5-dimethylphenoxy)ethoxy]phenyl}-6-[cyclopropyl-[2-(4-methoxyphenoxy)-ethyl]carbamoyl]-3,9-diazabicyclo[3.3.1]non-6-en-3-yl}-3-hydroxy-5-oxopentanoic acid formate salt

Synthesized according to typical procedures K, E and L from bicyclononene AL61 and 3-(*tert*-butyldimethylsilyloxy)glutaric anhydride. LC-MS: R_t = 0.86 ; ES+: 762.46

Example 501

1:1-Mixture of (*rac.*)-(3*R**)-5-{(*IR**, *5S**)-7-{4-[2-(2-chloro-4,5-dimethylphenoxy)ethoxy]phenyl}-6-(cyclopropylphenethylcarbamoyl)-3,9-diazabicyclo[3.3.1]non-6-en-3-yl}-3-hydroxy-5-oxopentanoic acid formate salt and (*rac.*)-(3*S**)-5-{(*IR**, *5S**)-7-{4-[2-(2-chloro-4,5-dimethylphenoxy)ethoxy]-

phenyl}-6-(cyclopropylphenethylcarbamoyl)-3,9-diazabicyclo[3.3.1]non-6-en-3-yl]-3-hydroxy-5-oxopentanoic acid formate salt

Synthesized according to typical procedures K, E and L from bicyclononene
5 AL64 and 3-(*tert*-butyldimethylsilyloxy)glutaric anhydride. LC-MS: R_t = 0.86 ;
ES+: 716.46

Example 502

10 1:1-Mixture of (*rac.*)-(3*R*^{*})-5-((1*R*^{*}, 5*S*^{*})-7-{4-[2-(2-chloro-4,5-dimethylphenoxy)ethoxy]phenyl}-6-{[2-(2-chlorophenyl)ethyl]cyclopropylcarbamoyl}-3,9-diazabicyclo[3.3.1]non-6-en-3-yl)-3-hydroxy-5-oxopentanoic acid formate salt and (*rac.*)-(3*R*^{*})-5-((1*R*^{*}, 5*S*^{*})-7-{4-[2-(2-chloro-4,5-dimethylphenoxy)ethoxy]phenyl}-6-{[2-(2-chlorophenyl)ethyl]cyclopropylcarbamoyl}-3,9-diazabicyclo[3.3.1]non-6-en-3-yl)-3-hydroxy-5-oxopentanoic acid formate salt

15

Synthesized according to typical procedures K, E and L from bicyclononene AL65 and 3-(*tert*-butyldimethylsilyloxy)glutaric anhydride. LC-MS: R_t = 0.88 ;
ES+: 750.41

20

Example 503

1:1-Mixture of (*rac.*)-(3*R*^{*})-5-((1*R*^{*}, 5*S*^{*})-7-{4-[2-(2-Chloro-4,5-dimethylphenoxy)ethoxy]phenyl}-6-{cyclopropyl-[2-(2,3-difluorophenyl)ethyl]carbamoyl}-3,9-diazabicyclo[3.3.1]non-6-en-3-yl)-3-hydroxy-5-oxopentanoic acid formate salt and (*rac.*)-(3*S*^{*})-5-((1*R*^{*}, 5*S*^{*})-7-{4-[2-(2-chloro-4,5-dimethylphenoxy)ethoxy]phenyl}-6-{cyclopropyl-[2-(2,3-difluorophenyl)ethyl]carbamoyl}-3,9-diazabicyclo[3.3.1]non-6-en-3-yl)-3-hydroxy-5-oxopentanoic acid formate salt

25

30

Synthesized according to typical procedures K, E and L from bicyclononene AL66 and 3-(*tert*-butyldimethylsilyloxy)glutaric anhydride. LC-MS: R_t = 0.88 ; ES+: 752.45

5 **Example 504**

- 1:1-Mixture of (*rac.*)-(3*R**)-5-((1*R**, 5*S**)-7-{4-[2-(2-chloro-4,5-dimethylphenoxy)ethoxy]phenyl}-6-{cyclopropyl-[2-(4-fluorophenyl)ethyl]-carbamoyl}-3,9-diazabicyclo[3.3.1]non-6-en-3-yl)-3-hydroxy-5-oxopentanoic acid formate salt and (*rac.*)-(3*S**)-5-((1*R**, 5*S**)-7-{4-[2-(2-chloro-4,5-dimethylphenoxy)ethoxy]phenyl}-6-{cyclopropyl-[2-(4-fluorophenyl)ethyl]-carbamoyl}-3,9-diazabicyclo[3.3.1]non-6-en-3-yl)-3-hydroxy-5-oxopentanoic acid formate salt
- 15 Synthesized according to typical procedures K, E and L from bicyclononene AL67 and 3-(*tert*-butyldimethylsilyloxy)glutaric anhydride. LC-MS: R_t = 0.87 ; ES+: 734.47

Example 505

- 20 1:1-Mixture of (*rac.*)-(3*R**)-5-((1*R**, 5*S**)-6-[(2-chlorobenzyl)cyclopropyl-carbamoyl]-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo-[3.3.1]non-6-en-9-yl)-3-hydroxy-5-oxo-pentanoic acid formate salt and (*rac.*)-(3*S**)-5-((1*R**, 5*S**)-6-[(2-chlorobenzyl)cyclopropylcarbamoyl]-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo-[3.3.1]non-6-en-9-yl)-3-hydroxy-5-oxo-pentanoic acid formate salt

- 25 Synthesized according to typical procedures K, E and L from bicyclononene BC and 3-(*tert*-butyldimethylsilyloxy)glutaric anhydride. LC-MS: R_t = 0.88 ; ES+: 726.32

Example 506

(rac.)-5-((1R, 5S*)-6-[(2-Chlorobenzyl)cyclopropylcarbamoyl]-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-en-9-yl)-2,2-dimethyl-5-oxopentanoic acid formate salt*

5

Synthesized according to typical procedures K and E from bicyclononene BC and 2,2-dimethylglutaric anhydride. LC-MS: R_t = 0.95 ; ES+: 738.37

Example 507

10

(rac.)-4-((1R, 5S*)-6-[(2-Chlorobenzyl)cyclopropylcarbamoyl]-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-en-9-yl)-4-oxo-butyric acid*

15

Synthesized according to typical procedures K and E from bicyclononene BC and succinic anhydride. LC-MS: R_t = 0.90 ; ES+: 696.32

Example 508

20

(rac.)-(1R, 5S*)-7-{4-[3-(2,3,6-Trifluorophenoxy)propyl]phenyl}-3,9-diaza-bicyclo[3.3.1]non-6-ene-6,9-dicarboxylic acid 6-[(2-chlorobenzyl)cyclopropyl-amide] 9-dimethylamide formate salt*

25

Synthesized according to typical procedures A and E from bicyclononene BC and dimethylcarbamoyl chloride. LC-MS: R_t = 0.94 ; ES+: 667.38

Example 509

30

(rac.)-(1R, 5S*)-6-[(2-Chlorobenzyl)cyclopropylcarbamoyl]-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-9-carboxylic acid methyl ester formate salt*

Synthesized according to typical procedures A and E from bicyclononene BC and methyl chloroformate. LC-MS: R_t = 0.96 ; ES+: 654.34

Example 510

5

(*rac.*)-(1*R*^{*}, 5*S*^{*})-6-[(2-Chlorobenzyl)cyclopropylcarbamoyl]-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-9-carboxylic acid ethyl ester formate salt

10 Synthesized according to typical procedures A and E from bicyclononene BC and ethyl chloroformate. LC-MS: R_t = 0.98 ; ES+: 668.37

Example 511

15 (*rac.*)-3-[(1*R*^{*}, 5*S*^{*})-(6-[(2-Chlorobenzyl)cyclopropylcarbamoyl]-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-9-carbonyl)amino]propionic acid ethyl ester formate salt

Synthesized according to typical procedures A and E from bicyclononene BC and 20 ethyl 4-isocyanatopropionate. LC-MS: R_t = 0.95 ; ES+: 739.36

Example 512

1:1-Mixture of (*rac.*)-(3*R*^{*})-5-{((1*R*^{*}, 5*S*^{*})-7-{4-[2-(2-chloro-4,5-dimethylphenoxy)ethoxy]phenyl}-6-[cyclopropyl(2-*o*-tolylethyl)carbamoyl]-3,9-diazabicyclo[3.3.1]non-6-en-3-yl)-3-hydroxy-5-oxopentanoic acid and (*rac.*)-(3*S*^{*})-5-{((1*R*^{*}, 5*S*^{*})-7-{4-[2-(2-chloro-4,5-dimethylphenoxy)ethoxy]phenyl}-6-[cyclopropyl(2-*o*-tolylethyl)carbamoyl]-3,9-diazabicyclo[3.3.1]non-6-en-3-yl)-3-hydroxy-5-oxopentanoic acid

Synthesized according to typical procedures K, E and L from bicyclononene **AL68** and 3-(*tert*-butyldimethylsilyloxy)glutaric anhydride. LC-MS: R_t = 0.88 ; ES+: 730.49

5 **Example 513**

1:1-Mixture of (*rac.*)-(3*R**)-5-{(*IR**, *SS**)-7-{4-[2-(2-chloro-4,5-dimethylphenoxy)ethoxy]phenyl}-6-[cyclopropyl-(3,5-dimethoxybenzyl)carbamoyl]-3,9-diazabicyclo[3.3.1]non-6-en-3-yl}-3-hydroxy-5-oxopentanoic acid formate salt and (*rac.*)-(3*S**)-5-{(*IR**, *SS**)-7-{4-[2-(2-chloro-4,5-dimethylphenoxy)ethoxy]phenyl}-6-[cyclopropyl-(3,5-dimethoxybenzyl)carbamoyl]-3,9-diazabicyclo[3.3.1]non-6-en-3-yl}-3-hydroxy-5-oxopentanoic acid formate salt

Synthesized according to typical procedures K, E and L from bicyclononene **AL70** and 3-(*tert*-butyldimethylsilyloxy)glutaric anhydride. LC-MS: R_t = 0.87; ES+: 762.47

Example 514

20 1:1-Mixture of (*rac.*)-(3*R**)-5-{(*IR**, *SS**)-7-{4-[2-(2-chloro-4,5-dimethylphenoxy)ethoxy]phenyl}-6-[cyclopropyl(2-*p*-tolylethyl)carbamoyl]-3,9-diazabicyclo[3.3.1]non-6-en-3-yl}-3-hydroxy-5-oxopentanoic acid formate salt and (*rac.*)-(3*S**)-5-{(*IR**, *SS**)-7-{4-[2-(2-chloro-4,5-dimethylphenoxy)ethoxy]phenyl}-6-[cyclopropyl(2-*p*-tolylethyl)carbamoyl]-3,9-diazabicyclo[3.3.1]non-6-en-3-yl}-3-hydroxy-5-oxopentanoic acid formate salt

25

Synthesized according to typical procedures K, E and L from bicyclononene **AL71** and 3-(*tert*-butyldimethylsilyloxy)glutaric anhydride. LC-MS: R_t = 0.88 ; ES+: 730.49

30

Example 515

1:1-Mixture of (*1R*, *5S*)-7-{4-[2-(2-chloro-4,5-dimethylphenoxy)ethoxy]-phenyl}-3-((*2S*, *4R*)-4-hydroxypyrrolidine-2-carbonyl)-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid (2-chlorobenzyl)cyclopropylamide formate salt and (*1S*, *5R*)-7-{4-[2-(2-chloro-4,5-dimethylphenoxy)ethoxy]phenyl}-3-((*2S*, *4R*)-4-hydroxypyrrolidine-2-carbonyl)-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid (2-chlorobenzyl)cyclopropylamide formate salt

Synthesized according to typical procedures G, E and L from bicyclononene AL56 and BOC-L-hydroxyproline. LC-MS: R_t = 0.81 ; ES+: 791.37

10

Example 516

1:1-Mixture of (*1R*, *5S*)-7-{4-[2-(2-chloro-4,5-dimethylphenoxy)ethoxy]-phenyl}-3-((*2S*, *4R*)-4-hydroxypyrrolidine-2-carbonyl)-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid cyclopropyl-(2-methylbenzyl)amide formate salt and (*1S*, *5R*)-7-{4-[2-(2-chloro-4,5-dimethylphenoxy)ethoxy]-phenyl}-3-((*2S*, *4R*)-4-hydroxypyrrolidine-2-carbonyl)-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid cyclopropyl-(2-methylbenzyl)amide formate salt

20

Synthesized according to typical procedures G, E and L from bicyclononene AL60 and BOC-L-hydroxyproline. LC-MS: R_t = 0.81 ; ES+: 699.45

25

(*rac.*)-(*1R*^{*}, *5S*^{*})-3-Acetyl-7-{4-[2-(2,3,5-trimethylphenoxy)ethyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid cyclopropylphenethylamide formate salt

30

Synthesized according to typical procedures A and E from bicyclononene AL47 and acetyl chloride. LC-MS: R_t = 0.92 ; ES+: 592.52

Example 518

(*rac.*)-(1*R*^{*, 5*S*^{*})-3-Acetyl-7-{4-[2-(2,3,5-trimethylphenoxy)ethyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid [2-(2-chlorophenyl)ethyl]-5 cyclopropylamide formate salt}

Synthesized according to typical procedures A and E from bicyclononene **AL48** and acetyl chloride. LC-MS: R_t = 0.94 ; ES+: 626.51

10 **Example 519**

(*rac.*)-(1*R*^{*, 5*S*^{*})-3-Acetyl-7-{4-[2-(2,3,5-trimethylphenoxy)ethyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid cyclopropyl-[2-(2,3-difluorophenyl)ethyl]amide formate salt}

15 Synthesized according to typical procedures A and E from bicyclononene **AL49** and acetyl chloride. LC-MS: R_t = 0.93 ; ES+: 628.54

20 **Example 520**

(*rac.*)-(1*R*^{*, 5*S*^{*})-3-Acetyl-7-{4-[2-(2,3,5-trimethylphenoxy)ethyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid cyclopropyl-(3,5-dimethoxybenzyl)amide formate salt}

25 Synthesized according to typical procedures A and E from bicyclononene **AL53** and acetyl chloride. LC-MS: R_t = 0.92 ; ES+: 638.54

Example 521

30 (*rac.*)-(1*R*^{*, 5*S*^{*})-3-acetyl-7-{4-[2-(2,3,5-Trimethylphenoxy)ethyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid cyclopropyl-(2-p-tolylethyl)-amide formate salt}

Synthesized according to typical procedures A and E from bicyclononene **AL54** and acetyl chloride. LC-MS: R_t = 0.94 ; ES+: 606.54

5 **Example 522**

(*rac.*)-(IR*, 5S*)-3-Acetyl-7-{4-[2-(2,3,5-trimethylphenoxy)ethyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid cyclopropyl-[2-(2-hydroxyethyl)benzyl]amide formate salt

10

Synthesized according to typical procedures A and E from bicyclononene **AL55** and acetyl chloride. LC-MS: R_t = 0.87 ; ES+: 622.55

Example 523

15

(*rac.*)-(IR*, 5S*)-3-Acetyl-7-{4-[2-(2-chloro-4,5-dimethylphenoxy)ethoxy]-phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid cyclopropyl-[2-(2-hydroxyethyl)benzyl]amide formate salt

20

Synthesized according to typical procedures A and E from bicyclononene **AL72** and acetyl chloride. LC-MS: R_t = 0.85 ; ES+: 658.47

Example 524

25

(*rac.*)-5-((IR*, 5S*)-6-(Cyclopropylphenethylcarbamoyl)-7-{4-[2-(2,3,5-trimethylphenoxy)ethyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-en-3-yl)-5-oxopentanoic acid formate salt

30

Synthesized according to typical procedures K and E from bicyclononene **AL47** and glutaric anhydride. LC-MS: R_t = 0.90 ; ES+: 664.56

Example 525

(rac.)-5-((1R, 5S*)-6-[2-(2-Chlorophenyl)ethyl]cyclopropylcarbamoyl)-7-{4-[2-(2,3,5-trimethylphenoxy)ethyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-en-3-yl)-5-oxopentanoic acid formate salt*

5

Synthesized according to typical procedures K and E from bicyclononene AL48 and glutaric anhydride. LC-MS: R_t = 0.91 ; ES+: 698.45

Example 526

10

(rac.)-5-((1R, 5S*)-6-[Cyclopropyl-[2-(2,3-difluorophenyl)ethyl]carbamoyl)-7-{4-[2-(2,3,5-trimethylphenoxy)ethyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-en-3-yl)-5-oxopentanoic acid*

15

Synthesized according to typical procedures K and E from bicyclononene AL49 and glutaric anhydride. LC-MS: R_t = 0.90 ; ES+: 700.50

Example 527

20

(rac.)-5-((1R, 5S*)-6-[Cyclopropyl-[2-(4-fluorophenyl)ethyl]carbamoyl)-7-{4-[2-(2,3,5-trimethylphenoxy)ethyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-en-3-yl)-5-oxopentanoic acid formate salt*

25

Synthesized according to typical procedures K and E from bicyclononene AL50 and glutaric anhydride. LC-MS: R_t = 0.90 ; ES+: 682.54

Example 528

30

(rac.)-5-((1R, 5S*)-6-[Cyclopropyl-(2-o-tolyethyl)carbamoyl]-7-{4-[2-(2,3,5-trimethylphenoxy)ethyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-en-3-yl)-5-oxopentanoic acid*

Synthesized according to typical procedures K and E from bicyclononene **AL51** and glutaric anhydride. LC-MS: R_t = 0.91 ; ES+: 678.55

Example 529

5

1:1-Mixture of (rac.)-5-((1R*, 5S*)-6-[cyclopropyl-((2R*)-2-hydroxy-2-phenylethyl)carbamoyl]-7-{4-[2-(2,3,5-trimethylphenoxy)ethyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-en-3-yl)-5-oxopentanoic acid formate salt and (rac.)-5-((1R*, 5S*)-6-[cyclopropyl-((2S*)-2-hydroxy-2-phenylethyl)carbamoyl]-7-{4-[2-(2,3,5-trimethylphenoxy)ethyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-en-3-yl)-5-oxopentanoic acid formate salt

Synthesized according to typical procedures K and E from bicyclononene **AL52** and glutaric anhydride. LC-MS: R_t = 0.85 ; ES+: 654.50

15

Example 530

(rac.)-5-((1R*, 5S*)-6-[Cyclopropyl-(3,5-dimethoxybenzyl)carbamoyl]-7-{4-[2-(2,3,5-trimethylphenoxy)ethyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-en-3-yl)-5-oxopentanoic acid formate salt

Synthesized according to typical procedures K and E from bicyclononene **AL53** and glutaric anhydride. LC-MS: R_t = 0.879 ; ES+: 710.50

25 Example 531

(rac.)-5-((1R*, 5S*)-6-[Cyclopropyl-(2-p-tolylethyl)carbamoyl]-7-{4-[2-(2,3,5-trimethylphenoxy)ethyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-en-3-yl)-5-oxopentanoic acid formate salt

30

Synthesized according to typical procedures K and E from bicyclononene **AL54** and glutaric anhydride. LC-MS: R_t = 0.91; ES+: 678.56

Example 532

5 *(rac.)-5-((1R*, 5S*)-6-{Cyclopropyl-[2-(2-hydroxyethyl)benzyl]carbamoyl}-7-{4-[2-(2,3,5-trimethylphenoxy)ethyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-en-3-yl)-5-oxo-pentanoic acid*

Synthesized according to typical procedures K and E from bicyclononene **AL55** and glutaric anhydride. LC-MS: R_t = 0.85 ; ES+: 694.53

10

Example 533

15 *(rac.)-5-((1R*, 5S*)-7-{4-[2-(2-Chloro-4,5-dimethylphenoxy)ethoxy]phenyl}-6-{cyclopropyl-[2-(2-hydroxyethyl)benzyl]carbamoyl}-3,9-diazabicyclo[3.3.1]-non-6-en-3-yl)-5-oxopentanoic acid formate salt*

Synthesized according to typical procedures K and E from bicyclononene **AL72** and glutaric anhydride. LC-MS: R_t = 0.84 ; ES+: 730.46

20 **Example 534**

25 *(rac.)-5-((1R*, 5S*)-6-(Cyclopropylphenethylcarbamoyl)-7-{4-[2-(2,3,5-tri-methylphenoxy)ethyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-en-3-yl)-2,2-di-methyl-5-oxopentanoic acid formate salt*

Synthesized according to typical procedures K and E from bicyclononene **AL47** and 2,2-dimethylglutaric anhydride. LC-MS: R_t = 0.93; ES+: 692.55

Example 535

30

(rac.)-5-((1R, 5S*)-6-[{2-(2-Chlorophenyl)ethyl]cyclopropylcarbamoyl}-7-{4-[2-(2,3,5-trimethylphenoxy)ethyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-en-3-yl)-2,2-dimethyl-5-oxopentanoic acid formate salt*

- 5 Synthesized according to typical procedures K and E from bicyclononene **AL48** and 2,2-dimethylglutaric anhydride. LC-MS: R_t = 0.95 ; ES+: 726.48

Example 536

- 10 *(rac.)-5-((1R*, 5S*)-6-[Cyclopropyl-[2-(2,3-difluorophenyl)ethyl]carbamoyl]-7-{4-[2-(2,3,5-trimethylphenoxy)ethyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-en-3-yl)-2,2-dimethyl-5-oxopentanoic acid formate salt*

Synthesized according to typical procedures K and E from bicyclononene **AL49** and 2,2-dimethylglutaric anhydride. LC-MS: R_t = 0.94; ES+: 728.51

Example 537

- 20 *(rac.)-5-((1R*, 5S*)-6-[Cyclopropyl-[2-(4-fluorophenyl)ethyl]carbamoyl]-7-{4-[2-(2,3,5-trimethylphenoxy)ethyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-en-3-yl)-2,2-dimethyl-5-oxopentanoic acid formate salt*

Synthesized according to typical procedures K and E from bicyclononene **AL50** and 2,2-dimethylglutaric anhydride. LC-MS: R_t = 0.94; ES+: 710.51

25

Example 538

- 30 *(rac.)-5-((1R*, 5S*)-6-[Cyclopropyl-(2-o-tolyethyl)carbamoyl]-7-{4-[2-(2,3,5-trimethylphenoxy)ethyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-en-3-yl)-2,2-dimethyl-5-oxopentanoic acid formate salt*

Synthesized according to typical procedures K and E from bicyclononene **AL51** and 2,2-dimethylglutaric anhydride. LC-MS: R_t = 0.95 ; ES+: 706.54

Example 539

5

(*rac.*)-5-((*1R*^{*}, *5S*^{*})-6-[Cyclopropyl-(3,5-dimethoxybenzyl)carbamoyl]-7-{4-[2-(2,3,5-trimethylphenoxy)ethyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-en-3-yl)-2,2-dimethyl-5-oxopentanoic acid formate salt

10 Synthesized according to typical procedures K and E from bicyclononene **AL53** and 2,2-dimethylglutaric anhydride. LC-MS: R_t = 0.93 ; ES+: 738.55

Example 540

15 (*rac.*)-5-((*1R*^{*}, *5S*^{*})-6-[Cyclopropyl-(2-p-tolyethyl)carbamoyl]-7-{4-[2-(2,3,5-trimethylphenoxy)ethyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-en-3-yl)-2,2-dimethyl-5-oxopentanoic acid formate salt

20 Synthesized according to typical procedures K and E from bicyclononene **AL54** and 2,2-dimethylglutaric anhydride. LC-MS: R_t = 0.95 ; ES+: 706.53

Example 541

25 (*rac.*)-5-((*1R*^{*}, *5S*^{*})-7-{4-[2-(2-Chloro-4,5-dimethylphenoxy)ethoxy]phenyl}-6-[cyclopropyl-[2-(2-hydroxyethyl)benzyl]carbamoyl]-3,9-diazabicyclo[3.3.1]-non-6-en-3-yl)-2,2-dimethyl-5-oxopentanoic acid formate salt

Synthesized according to typical procedures K and E from bicyclononene **AL72** and 2,2-dimethylglutaric anhydride. LC-MS: R_t = 0.87 ; ES+: 758.46

30

Example 542

(*rac.*)-(1*R*^{*}, 5*S*^{*})-3-(4-Carbamoylbutyryl)-7-{4-[2-(2,3,5-trimethylphenoxy)-ethyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid cyclopropyl-phenethylamide formate salt

5 Synthesized according to typical procedures G and E from bicyclononene AL47 and 4-carbamoylbutyric acid. LC-MS: R_t = 0.88 ; ES+: 663.55

Example 543

10 **(*rac.*)-(1*R*^{*}, 5*S*^{*})-3-(4-Carbamoylbutyryl)-7-{4-[2-(2,3,5-trimethylphenoxy)-ethyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid [2-(2-chlorophenyl)ethyl]cyclopropylamide formate salt**

15 Synthesized according to typical procedures G and E from bicyclononene AL48 and 4-carbamoylbutyric acid. LC-MS: R_t = 0.90 ; ES+: 697.49

Example 544

20 **(*rac.*)-(1*R*^{*}, 5*S*^{*})-3-(4-Carbamoylbutyryl)-7-{4-[2-(2,3,5-trimethylphenoxy)-ethyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid cyclopropyl-[2-(2,3-difluorophenyl)ethyl]amide formate salt**

25 Synthesized according to typical procedures G and E from bicyclononene AL49 and 4-carbamoylbutyric acid. LC-MS: R_t = 0.89 ; ES+: 699.49

Example 545

30 **(*rac.*)-(1*R*^{*}, 5*S*^{*})-3-(4-Carbamoylbutyryl)-7-{4-[2-(2,3,5-trimethylphenoxy)-ethyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid cyclopropyl-[2-(4-fluorophenyl)ethyl]amide formate salt**

Synthesized according to typical procedures G and E from bicyclononene **AL50** and 4-carbamoylbutyric acid. LC-MS: R_t = 0.88 ; ES+: 681.54

Example 546

5

(*rac.*)-(1*R*^{*}, 5*S*^{*})-3-(4-Carbamoylbutyryl)-7-{4-[2-(2,3,5-trimethylphenoxy)-ethyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid cyclopropyl-(3,5-dimethoxybenzyl)amide formate salt

10 Synthesized according to typical procedures G and E from bicyclononene **AL53** and 4-carbamoylbutyric acid. LC-MS: R_t = 0.87 ; ES+: 709.50

Example 547

15 (*rac.*)-(1*R*^{*}, 5*S*^{*})-3-(4-Carbamoylbutyryl)-7-{4-[2-(2,3,5-trimethylphenoxy)-ethyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid cyclopropyl-(2-*p*-tolylethyl)amide formate salt

20 Synthesized according to typical procedures G and E from bicyclononene **AL54** and 4-carbamoylbutyric acid. LC-MS: R_t = 0.89 ; ES+: 677.57

Example 548

25 (*rac.*)-(1*R*^{*}, 5*S*^{*})-3-(4-Carbamoylbutyryl)-7-{4-[2-(2,3,5-trimethylphenoxy)-ethyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid cyclopropyl-[2-(2-hydroxyethyl)benzyl]amide formate salt

Synthesized according to typical procedures G and E from bicyclononene **AL55** and 4-carbamoylbutyric acid. LC-MS: R_t = 0.83 ; ES+: 693.53

30

Example 549

(rac.)-(1R, 5S*)-3-(4-Carbamoylbutyryl)-7-{4-[2-(2-chloro-4,5-dimethyl-phenoxy)ethoxy]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid cyclopropyl-[2-(2-hydroxyethyl)benzyl]amide formate salt*

5 Synthesized according to typical procedures G and E from bicyclononene AL72 and 4-carbamoylbutyric acid. LC-MS: R_t = 0.82 ; ES+: 729.48

Example 550

10 *(rac.)-Acetic acid (1R*, 5S*)-2-(2-[(3-acetyl-7-{4-[2-(2,3,5-trimethyl-phenoxy)ethyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carbonyl)cyclo-propylamino]methyl]phenyl)ethyl ester formate salt*

15 Synthesized according to typical procedures A and E from bicyclononene AL and acetyl chloride. LC-MS: R_t = 0.93 ; ES+: 664.55

Example 551

20 *(rac.)-Acetic acid (1R*, 5S*)-2-(2-[(3-acetyl-7-{4-[2-(2-chloro-4,5-dimethyl-phenoxy)ethoxy]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carbonyl)cyclo-propylamino]methyl]phenyl)ethyl ester formate salt*

25 Synthesized according to typical procedures A and E from bicyclononene AL72 and acetyl chloride. LC-MS: R_t = 0.91 ; ES+: 700.44

Example 552

30 *(rac.)-(1R*, 5S*)-7-{4-[2-(2,3,5-Trimethylphenoxy)ethyl]phenyl}-3,9-diaza-bicyclo[3.3.1]non-6-ene-6-carboxylic acid cyclopropyl-[2-(4-fluorophenyl)-ethyl]amide formate salt*

Synthesized according to typical procedure E from bicyclononene **AK50**. LC-MS: R_t = 0.86 ; ES+: 568.50

Example 553

5

(*rac.*)-(1*R*^{*}, 5*S*^{*})-7-{4-[2-(2,3,5-Trimethylphenoxy)ethyl]phenyl}-3,9-diaza-bicyclo[3.3.1]non-6-ene-6-carboxylic acid cyclopropyl-(3,5-dimethoxybenzyl)-amide formate salt

10 Synthesized according to typical procedure E from bicyclononene **AK53**. LC-MS: R_t = 0.85 ; ES+: 596.53

Example 554

15 (*rac.*)-Acetic acid (1*R*^{*}, 5*S*^{*})-2-(2-[(3-acetyl-7-{4-[2-(4-bromophenoxy)-ethoxy]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carbonyl)cyclopropyl-amino]methyl)phenyl)ethyl ester formate salt

20 Synthesized according to typical procedures A and E from bicyclononene **AL87** and acetyl chloride. LC-MS: R_t = 0.89 ; ES+: 716.35

Example 555

25 1:1 Mixture of (*rac.*)-5-{(1*R*^{*}, 5*S*^{*})-7-{4-[2-(2-chloro-4,5-dimethylphenoxy)-ethoxy]phenyl}-6-[((2*R*^{*})-2-hydroxy-2-phenylethyl)methylcarbamoyl]-3,9-diaza-bicyclo[3.3.1]non-6-en-3-yl}-2,2-dimethyl-5-oxopentanoic acid formate salt and (*rac.*)-5-{(1*R*^{*}, 5*S*^{*})-7-{4-[2-(2-chloro-4,5-dimethylphenoxy)-ethoxy]phenyl}-6-[((2*S*^{*})-2-hydroxy-2-phenylethyl)methylcarbamoyl]-3,9-diaza-bicyclo[3.3.1]non-6-en-3-yl}-2,2-dimethyl-5-oxopentanoic acid formate salt

Synthesized according to typical procedures K and E from bicyclononene **AL69** and 2,2-dimethylglutaric anhydride. LC-MS: R_t = 0.88 ; ES+: 718.47

Example 556

5

(*rac.*)-(IR*, 5S*)-3-Acetyl-7-{4-[3-(2-bromo-5-fluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid (2-chlorobenzyl)cyclopropylamide formate salt

10 Synthesized according to typical procedures G and E from bicyclononene **AL20** and acetyl chloride. LC-MS: R_t = 3.63; ES+: 680.28.

Example 557

15 (*rac.*)-(IR*, 5S*)-7-{4-[3-(2,3,6-Trifluorophenoxy)propyl]phenyl}-3,9-diaza-bicyclo[3.3.1]non-6-ene-6-carboxylic acid benzylcyclopropylamide formate salt

Synthesized according to typical procedure E from bicyclononene **AK3**. LC-MS:
20 R_t = 0.83 ; ES+: 562.38

Example 558

(*rac.*)-(IR*, 5S*)-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diaza-bicyclo[3.3.1]non-6-ene-6-carboxylic acid (2-chlorobenzyl)ethylamide formate salt

Synthesized according to typical procedure E from bicyclononene **AK4**. LC-MS:
R_t = 0.84 ; ES+: 584.35

30

Example 559

(*rac.*)-(1*R*^{*}, 5*S*^{*})-7-{4-[3-(2,3,6-Trifluorophenoxy)propyl]phenyl}-3,9-diaza-bicyclo[3.3.1]non-6-ene-6-carboxylic acid cyclopropyl-(2-fluorobenzyl)amide formate salt

5 Synthesized according to typical procedure E from bicyclononene AK5. LC-MS:
R_t = 0.83 ; ES+: 580.38

Example 560

10 **(*rac.*)-(1*R*^{*}, 5*S*^{*})-7-{4-[3-(2,3,6-Trifluorophenoxy)propyl]phenyl}-3,9-diaza-bicyclo[3.3.1]non-6-ene-6-carboxylic acid cyclopropyl-(3-trifluoromethylbenzyl)amide**

15 Synthesized according to typical procedure E from bicyclononene AK6. LC-MS:
R_t = 0.86 ; ES+: 630.43

Example 561

20 **(*rac.*)-(1*R*^{*}, 5*S*^{*})-7-{4-[3-(2,3,6-Trifluorophenoxy)propyl]phenyl}-3,9-diaza-bicyclo[3.3.1]non-6-ene-6-carboxylic acid cyclopropyl-(2-methylbenzyl)amide formate salt**

25 Synthesized according to typical procedure E from bicyclononene AK7. LC-MS:
R_t = 0.84 ; ES+: 576.42

25

Example 562

30 **(*rac.*)-(1*R*^{*}, 5*S*^{*})-7-{4-[3-(2,3,6-Trifluorophenoxy)propyl]phenyl}-3,9-diaza-bicyclo[3.3.1]non-6-ene-6-carboxylic acid cyclopropyl-[2-(4-methoxyphenoxy)ethyl]amide formate salt**

Synthesized according to typical procedure E from bicyclononene **AK8**. LC-MS:
R_t = 0.83 ; ES+: 622.45

Example 563

5

(*rac.*)-(1*R*^{*}, 5*S*^{*})-7-{4-[3-(2,3,6-Trifluorophenoxy)propyl]phenyl}-3,9-diaza-bicyclo[3.3.1]non-6-ene-6-carboxylic acid [2-(2-chlorophenyl)ethyl]-cyclopropylamide formate salt

10 Synthesized according to typical procedure E from bicyclononene **AK13**. LC-MS: R_t = 0.86 ; ES+: 610.39

Example 564

15 (*rac.*)-(1*R*^{*}, 5*S*^{*})-7-{4-[3-(2,3,6-Trifluorophenoxy)propyl]phenyl}-3,9-diaza-bicyclo[3.3.1]non-6-ene-6-carboxylic acid cyclopropyl-[2-(2,3-difluorophenyl)ethyl]amide formate salt

20 Synthesized according to typical procedure E from bicyclononene **AK14**. LC-MS: R_t = 0.85 ; ES+: 612.44

Example 565

25 (*rac.*)-(1*R*^{*}, 5*S*^{*})-7-{4-[3-(2,3,6-Trifluorophenoxy)propyl]phenyl}-3,9-diaza-bicyclo[3.3.1]non-6-ene-6-carboxylic acid cyclopropyl-(2-*o*-tolylethyl)amide formate salt

Synthesized according to typical procedure E from bicyclononene **AK16**. LC-MS: R_t = 0.85 ; ES+: 590.42

30

Example 566

(*rac.*)-(1*R*^{*}, 5*S*^{*})-7-{4-[3-(2,3,6-Trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid cyclopropyl-(2-*p*-tolylethyl)amide formate salt

5 Synthesized according to typical procedures G and E from bicyclononene **AK18**.
LC-MS: R_t = 0.85 ; ES+: 590.42

Example 567

10 (*rac.*)-(1*R*^{*}, 5*S*^{*})-3-Acetyl-7-{4-[3-(2,5-difluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid (2-chlorobenzyl)cyclopropylamide formate salt

15 Synthesized according to typical procedures F and E from bicyclononene **BB1** and 2,5-difluorophenol. LC-MS: R_t = 0.89 ; ES+: 614.40

Example 568

20 (*rac.*)-(1*R*^{*}, 5*S*^{*})-3-Acetyl-7-{4-[3-(2,3-dichlorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid (2-chlorobenzyl)cyclopropylamide formate salt

25 Synthesized according to typical procedures F and E from bicyclononene **BB1** and 2,3-dichlorophenol. LC-MS: R_t = 0.94 ; ES+: 654.32

25

Example 569

30 (*rac.*)-(1*R*^{*}, 5*S*^{*})-3-Acetyl-7-{4-[3-(2-chloro-5-fluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid (2-chlorobenzyl)cyclopropylamide formate salt

Synthesized according to typical procedures F and E from bicyclononene **BB1** and 2-chloro-5-fluorophenol. LC-MS: R_t = 0.93 ; ES+: 636.36

Example 570

5

(*rac.*)-(1*R*^{*}, 5*S*^{*})-3-Acetyl-7-{4-[3-(3-cyanopyridin-2-yloxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid (2-chlorobenzyl)cyclopropylamide formate salt

10 Synthesized according to typical procedures F and E from bicyclononene **BB1** and 2-hydroxynicotinonitrile. LC-MS: R_t = 0.86 ; ES+: 610.42

Example 571

15 (*rac.*)-(1*R*^{*}, 5*S*^{*})-7-{4-[3-(2,3,6-Trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-3,6-dicarboxylic acid 6-[(2-chlorobenzyl)cyclopropylamide] 3-dimethylamide formate salt

20 Synthesized according to typical procedures A and E from bicyclononene **AL2** and dimethylcarbamoyl chloride. LC-MS: R_t = 0.91 ; ES+: 667.42

Example 572

25 (*rac.*)-(1*R*^{*}, 5*S*^{*})-7-{4-[3-(2,3,6-Trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-3,6-dicarboxylic acid 6-[(2-chlorobenzyl)cyclopropylamide] 3-diethylamide formate salt

Synthesized according to typical procedures A and E from bicyclononene **AL2** and diethylcarbamoyl chloride. LC-MS: R_t = 0.95 ; ES+: 695.44

30

Example 573

(rac.)-(1*R*^{*, 5*S*^{*})-6-[(2-Chlorobenzyl)cyclopropylcarbamoyl]-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-3-carboxylic acid methyl ester formate salt}

5 Synthesized according to typical procedures A and E from bicyclononene AL2 and methyl chloroformate. LC-MS: R_t = 0.91 ; ES+: 654.37

Example 574

10 **(rac.)-(1*R*^{, 5*S*^{*})-6-[(2-Chlorobenzyl)cyclopropylcarbamoyl]-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-3-carboxylic acid ethyl ester formate salt}**

15 Synthesized according to typical procedures A and E from bicyclononene AL2 and ethyl chloroformate. LC-MS: R_t = 0.93 ; ES+: 668.40

Example 575

20 **(rac.)-(1*R*^{, 5*S*^{*})-3-Methanesulfonyl-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]-phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid (2-chlorobenzyl)-cyclopropylamide formate salt}**

25 Synthesized according to typical procedures A and E from bicyclononene AL2 and methanesulfonyl chloride. LC-MS: R_t = 0.92 ; ES+: 674.37

25

Example 576

30 **(rac.)-(1*R*^{, 5*S*^{*})-3-Ethanesulfonyl-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]-phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid (2-chlorobenzyl)-cyclopropylamide formate salt}**

Synthesized according to typical procedures A and E from bicyclononene **AL2** and ethylsulfonyl chloride. LC-MS: R_t = 0.93 ; ES+: 688.36

Example 577

5

(*rac.*)-5-((*1R*^{*,} *5S*^{*)}-6-[(2-Chlorobenzyl)cyclopropylcarbamoyl]-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-en-3-yl)-5-oxo-pentanoic acid ethyl ester formate salt

10 Synthesized according to typical procedures A and E from bicyclononene **AL2** and glutaric acid monoethylester chloride. LC-MS: R_t = 0.94 ; ES+: 738.41

Example 578

15 (*rac.*)-4-((*1R*^{*,} *5S*^{*)}-6-[(2-Chlorobenzyl)cyclopropylcarbamoyl]-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-en-3-yl)-4-oxo-butyric acid formate salt

20 Synthesized according to typical procedures K and E from bicyclononene **AL2** and succinic anhydride. LC-MS: R_t = 0.87 ; ES+: 696.36

Example 579

25 (*rac.*)-3-[((*1R*^{*,} *5S*^{*)}-6-[(2-Chlorobenzyl)cyclopropylcarbamoyl]-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-3-carbonyl)amino]propionic acid ethyl ester formate salt

30 Synthesized according to typical procedures A and E from bicyclononene **AL2** and ethyl 4-isocyanatopropionate. LC-MS: R_t = 0.93 ; ES+: 739.41

Example 580

(*rac.*)4-[(*1R*^{*}, *5S*^{*})-6-[(2-Chlorobenzyl)cyclopropylcarbamoyl]-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-3-carbonyl]amino]butyric acid ethyl ester formate salt

5 Synthesized according to typical procedures A and E from bicyclononene **AL2** and ethyl 4-isocyanatobutyrate. LC-MS: R_t = 0.93 ; ES+: 753.39

Example 581

10 **(*rac.*)-(1*R*^{*}, 5*S*^{*})-3-(3-Carbamoylpropionyl)-7-{4-[3-(2,3,6-trifluorophenoxy)-propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid (2-chlorobenzyl)cyclopropylamide formate salt**

15 Synthesized according to typical procedures G and E from bicyclononene **AL2** and succinamic acid. LC-MS: R_t = 0.85 ; ES+: 695.38

Example 582

20 **(*rac.*)-(1*R*^{*}, 5*S*^{*})-3-(2-Hydroxyacetyl)-7-{4-[3-(2,3,6-trifluorophenoxy)-propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid (2-chlorobenzyl)cyclopropylamide formate salt**

25 Synthesized according to typical procedures G and E from bicyclononene **AL2** and glycolic acid. LC-MS: R_t = 0.88 ; ES+: 654.36

25

Example 583

30 **(*1S*, 5*R*)-3-((3*R*)-3-Hydroxybutyryl)-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid (2-chlorobenzyl)-cyclopropylamide**

Synthesized according to typical procedures G and E from bicyclononene **AL2** and (*3R*)-3-hydroxybutyric acid. LC-MS: R_t = 0.88; ES+: 682.40.

Example 584

5

1:1-Mixture of (*rac.*)-(1*R*^{*}, 5*S*^{*})-3-((1*R*^{*}, 2*S*^{*})-2-hydroxycyclopentane-carbonyl)-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid (2-chlorobenzyl)cyclopropylamide formate salt and (*rac.*)-(1*R*^{*}, 5*S*^{*})-3-((1*S*^{*}, 2*R*^{*})-2-hydroxycyclopentanecarbonyl)-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid (2-chlorobenzyl)cyclopropylamide formate salt

10

15

Synthesized according to typical procedures G and E from bicyclononene **AL2** and *cis*-2-hydroxy-1-cyclopentane carboxylic acid. LC-MS: R_t = 0.91; ES+:

708.39.

Example 585

20

(*rac.*)-(1*R*^{*}, 5*S*^{*})-9-Acetyl-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid (2-chlorobenzyl)cyclopropylamide formate salt

25

Synthesized according to typical procedures A and E from bicyclononene **BC** and acetyl chloride. LC-MS: R_t = 0.92; ES+: 638.34.

Example 586

30

(*rac.*)-5-((1*R*^{*}, 5*S*^{*})-6-[(2-Chlorobenzyl)cyclopropylcarbamoyl]-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-en-9-yl)-5-oxo-pentanoic acid formate salt

Synthesized according to typical procedures K and E from bicyclononene BC and glutaric anhydride. LC-MS: R_t = 0.91; ES+: 710.33.

Example 587

5

(rac.)-5-((1R, 5S*)-6-[(2-Chlorobenzyl)cyclopropylcarbamoyl]-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-en-9-yl)-5-oxo-pentanoic acid methyl ester formate salt*

10 Synthesized according to typical procedures A and E from bicyclononene BC and glutaric acid monomethylester chloride. LC-MS: R_t = 0.95; ES+: 724.33.

Example 588

15

(rac.)-5-((1R, 5S*)-6-[(2-Chlorobenzyl)cyclopropylcarbamoyl]-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-en-9-yl)-5-oxo-pentanoic acid ethyl ester formate salt*

20

Synthesized according to typical procedures A and E from bicyclononene BC and glutaric acid monoethylester chloride. LC-MS: R_t = 0.97; ES+: 738.38.

Example 589

25

(rac.)-4-[((1R, 5S*)-6-[(2-Chlorobenzyl)cyclopropylcarbamoyl]-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-9-carbonyl)-amino]butyric acid ethyl ester*

30

Synthesized according to typical procedures A and E from bicyclononene BC and ethyl 4-isocyanatobutyrate. LC-MS: R_t = 0.95; ES+: 753.37.

Example 590

(*rac.*)-(1*R*^{*}, 5*S*^{*})-3-Formyl-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid (2-chlorobenzyl)cyclopropylamide

5 Synthesized according to typical procedure E from bicyclononene **AK2**. Obtained as side-product after purification by HPLC. LC-MS: R_t = 0.89; ES+: 624.36.

Example 591

10 **(*rac.*)-3-[((1*R*^{*}, 5*S*^{*})-6-[(2-Chlorobenzyl)cyclopropylcarbamoyl]-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-3-carbonyl)amino]propionic acid formate salt**

15 Synthesized according to typical procedures M and E from **Example 579**, then typical procedure E. LC-MS: R_t = 0.87; ES+: 711.31.

Example 592

20 **(*rac.*)-3-[((1*R*^{*}, 5*S*^{*})-6-[(2-Chlorobenzyl)cyclopropylcarbamoyl]-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-9-carbonyl)amino]propionic acid formate salt**

Synthesized according to typical procedures G, M and E from bicyclononene **BC** and ethyl 4-isocyanatopropionate. LC-MS: R_t = 0.88; ES+: 711.33.

25

Example 593

30 **(*rac.*)-4-[((1*R*^{*}, 5*S*^{*})-6-[(2-Chlorobenzyl)cyclopropylcarbamoyl]-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-9-carbonyl)amino]butyric acid formate salt**

Synthesized according to typical procedures M and E from **Example 580**. LC-MS: R_t = 0.89; ES+: 725.35.

Example 594

5

(rac.)-(IR, 5S*)-3-acetyl-7-{4-[2-(2,3,6-Trifluorophenoxy)ethyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid (2-chlorobenzyl)cyclopropylamide formate salt*

10 Synthesized according to typical procedures F and E from bicyclononene **BB2** and 2,3,6-trifluorophenol. LC-MS: R_t = 0.89; ES+: 624.37.

Example 595

15 *(rac.)-(IR*, 5S*)-3-Acetyl-7-{4-[2-(2,3-dimethylphenoxy)ethyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid (2-chlorobenzyl)cyclopropylamide formate salt*

20 Synthesized according to typical procedures F and E from bicyclononene **BB2** and 2,3-dimethylphenol. LC-MS: R_t = 0.91; ES+: 598.42.

Example 596

25 *(rac.)-(IR*, 5S*)-3-Acetyl-7-{4-[2-(2,5-dimethylphenoxy)ethyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid (2-chlorobenzyl)cyclopropylamide*

Synthesized according to typical procedures F and E from bicyclononene **BB2** and 2,5-dimethylphenol. LC-MS: R_t = 0.91; ES+: 598.43.

30

Example 597

(*rac.*)-(1*R*^{*,}, 5*S*^{*})-3-Acetyl-7-{4-[2-(2-chloro-4,5-dimethylphenoxy)ethyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid (2-chlorobenzyl)cyclopropylamide formate salt

5 Synthesized according to typical procedures G and E from bicyclononene BB2 and 2-chloro-4,5-dimethylphenol. LC-MS: R_t = 0.93; ES+: 632.39.

Example 598

10 **(*rac.*)-(1*R*^{*,}, 5*S*^{*})-3-Acetyl-7-{4-[2-(2,4-dichlorophenoxy)ethyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid (2-chlorobenzyl)cyclopropylamide formate salt**

15 Synthesized according to typical procedures F and E from bicyclononene BB2 and 2,4-dichlorophenol. LC-MS: R_t = 0.91; ES+: 640.32.

Example 599

20 **(*rac.*)-(1*R*^{*,}, 5*S*^{*})-3-Acetyl-7-{4-[2-(2,3-dichlorophenoxy)ethyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid (2-chlorobenzyl)cyclopropylamide formate salt**

25 Synthesized according to typical procedures F and E from bicyclononene BB2 and 2,3-dichlorophenol. LC-MS: R_t = 0.91; ES+: 640.34.

Example 600

30 **(*rac.*)-(1*R*^{*,}, 5*S*^{*})-3-Acetyl-7-{4-[2-(2,6-difluorophenoxy)ethyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid (2-chlorobenzyl)cyclopropylamide formate salt**

Synthesized according to typical procedures F and E from bicyclononene BB2 and 2,6-difluorophenol. LC-MS: R_t = 0.88; ES+: 606.39.

Example 601

5

(*rac.*)-(IR*, 5S*)-3-Acetyl-7-{4-[2-(2,5-difluorophenoxy)ethyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid (2-chlorobenzyl)cyclopropylamide formate salt

10 Synthesized according to typical procedures F and E from bicyclononene BB2 and 2,5-difluorophenol. LC-MS: R_t = 0.88; ES+: 606.40.

Example 602

15 (*rac.*)-(IR*, 5S*)-3-Acetyl-7-{4-[2-(3,5-dichlorophenoxy)ethyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid (2-chlorobenzyl)cyclopropylamide

20 Synthesized according to typical procedures F and E from bicyclononene BB2 and 3,5-dichlorophenol. LC-MS: R_t = 0.93; ES+: 638.32.

Example 603

25 (*rac.*)-(IR*, 5S*)-3-Acetyl-7-{4-[2-(2-chloro-5-methylphenoxy)ethyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid (2-chlorobenzyl)cyclopropylamide formate salt

Synthesized according to typical procedures F and E from bicyclononene BB2 and 2-chloro-5-methylphenol. LC-MS: R_t = 0.91; ES+: 618.40.

30

Example 604

(rac.)-(1R, 5S*)-3-Acetyl-7-{4-[2-(2-chloro-5-fluorophenoxy)ethyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid (2-chlorobenzyl)cyclopropylamide formate salt*

- 5 Synthesized according to typical procedures F and E from bicyclononene **BB2** and 2-chloro-5-fluorophenol. LC-MS: R_t = 0.89; ES+: 622.33.

Example 605

- 10 *(rac.)-(1R*, 5S*)-3-Acetyl-7-{4-[2-(2,3,6-trichlorophenoxy)ethyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid (2-chlorobenzyl)cyclopropylamide formate salt*

- 15 Synthesized according to typical procedures F and E from bicyclononene **BB2** and 2,5,6-trichlorophenol. LC-MS: R_t = 0.93; ES+: 674.27.

Example 606

- 20 *(rac.)-(1R*, 5S*)-9-(4-Carbamoylbutyryl)-7-{4-[2-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid (2-chlorobenzyl)cyclopropylamide formate salt*

- Synthesized according to typical procedures G and E from bicyclononene **BC** and 4-carbamoylbutyric acid. LC-MS: R_t = 0.88; ES+: 709.39.

25

Example 607

- 30 *(rac.)-(1R*, 5S*)-9-(3-Carbamoylpropionyl)-7-{4-[2-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid (2-chlorobenzyl)cyclopropylamide formate salt*

Synthesized according to typical procedures G and E from bicyclononene BC and succinamic acid. LC-MS: R_t = 0.88; ES+: 695.36.

Example 608

5

(*rac.*)-(1*R*^{*}, 5*S*^{*})-9-(2-Hydroxyacetyl)-7-{4-[2-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid (2-chlorobenzyl)cyclopropylamide formate salt

10 Synthesized according to typical procedures G and E from bicyclononene BC and glycolic acid. LC-MS: R_t = 0.90; ES+: 654.37.

Example 609

15 (1*R*, 5*S*)-9-((3*S*)-3-Hydroxybutyryl)-7-{4-[2-(2,3,6-trifluorophenoxy)propyl]-phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid (2-chlorobenzyl)-cyclopropylamide formate salt

20 Synthesized according to typical procedures G and E from bicyclononene BC and (3*R*)-3-hydroxybutyric acid. LC-MS: R_t = 0.91; ES+: 682.41.

Example 610

25 (1*R*, 5*S*)-9-Methanesulfonyl-7-{4-[2-(2,3,6-trifluorophenoxy)propyl]-phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid (2-chlorobenzyl)-cyclopropylamide formate salt

Synthesized according to typical procedures A and E from bicyclononene BC and methylsulfonyl chloride. LC-MS: R_t = 0.95; ES+: 674.38.

30

Example 611

(rac.)-(IR, 5S*)-9-Ethanesulfonyl-7-{4-[2-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid (2-chlorobenzyl)-cyclopropylamide formate salt*

- 5 Synthesized according to typical procedures A and E from bicyclononene BC and ethylsulfonyl chloride. LC-MS: R_t = 0.97; ES+: 688.37.

Example 612

- 10 *(rac.)-(IR*, 5S*)-3-Acetyl-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid cyclopropyl-(2-fluoro-5-methoxybenzyl)amide formate salt*

Synthesized according to typical procedures H and E from bicyclononene AJ2
15 and cyclopropyl(2-fluoro-5-methoxybenzyl)amine. LC-MS: R_t = 0.89; ES+: 652.31.

Example 613

- 20 *(rac.)-(IR*, 5S*)-3-Acetyl-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid cyclopropyl-(3-methoxybenzyl)amide formate salt*

Synthesized according to typical procedures H and E from bicyclononene AJ2
25 and cyclopropyl(3-methoxybenzyl)amine. LC-MS: R_t = 0.88; ES+: 634.34.

Example 614

- 30 *(rac.)-(IR*, 5S*)-3-Acetyl-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid cyclopropyl-(2-methoxybenzyl)amide formate salt*

Synthesized according to typical procedures H and E from bicyclononene **AJ2** and cyclopropyl(2-methoxybenzyl)amine. LC-MS: $R_t = 0.89$; ES+: 634.32.

Example 615

5

(*rac.*)-(1*R**, 5*S**)-3-Acetyl-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid cyclopropyl-(3,4-dimethoxybenzyl)amide formate salt

10 Synthesized according to typical procedures H and E from bicyclononene **AJ2** and cyclopropyl(3,4-dimethoxybenzyl)amine. LC-MS: $R_t = 0.86$; ES+: 664.34.

Example 616

15 (*rac.*)-(1*R**, 5*S**)-3-Acetyl-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid (2-chloro-3-trifluoromethylbenzyl)cyclopropylamide formate salt

20 Synthesized according to typical procedures H and E from bicyclononene **AJ2** and (2-chloro-3-trifluoromethylbenzyl)cyclopropylamine. LC-MS: $R_t = 0.94$; ES+: 706.20.

Example 617

25 (*rac.*)-(1*R**, 5*S**)-3-Acetyl-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid benzo[1,3]dioxol-5-ylmethylcyclopropylamide formate salt

30 Synthesized according to typical procedures H and E from bicyclononene **AJ2** and (6-chlorobenzo[1,3]dioxol-5-ylmethyl)cyclopropylamine. LC-MS: $R_t = 0.91$; ES+: 682.28.

Example 618

(*rac.*)-(1*R*^{*}, 5*S*^{*})-3-Acetyl-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid cyclopropyl-(5-fluoro-2-methoxybenzyl)amide formate salt

Synthesized according to typical procedures H and E from bicyclononene **AJ2** and cyclopropyl(2-methoxy-5-fluorobenzyl)amine. LC-MS: R_t = 0.90; ES+: 652.32.

10

Example 619

(*rac.*)-(1*R*^{*}, 5*S*^{*})-3-Acetyl-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid (2-chloro-6-fluorobenzyl)-cyclopropylamide formate salt

Synthesized according to typical procedures H and E from bicyclononene **AJ2** and cyclopropyl(2-chloro-6-fluorobenzyl)amine. LC-MS: R_t = 0.89; ES+: 656.30.

20 **Example 620**

(*rac.*)-(1*R*^{*}, 5*S*^{*})-3-Acetyl-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid (2-bromobenzyl)cyclopropyl-amide formate salt

25

Synthesized according to typical procedures H and E from bicyclononene **AJ2** and cyclopropyl(2-bromobenzyl)amine. LC-MS: R_t = 0.91; ES+: 684.23.

Example 621

30

(*rac.*)-(1*R*^{*}, 5*S*^{*})-3-Acetyl-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid cyclopropyl-(2,6-difluorobenzyl)amide formate salt

5 Synthesized according to typical procedures H and E from bicyclononene AJ2 and cyclopropyl(2,6-difluorobenzyl)amine. LC-MS: R_t = 0.88; ES+: 640.29.

Example 622

10 **(*rac.*)-(1*R*^{*}, 5*S*^{*})-3-Acetyl-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid cyclopropyl-(2,3-dimethylbenzyl)amide formate salt**

15 Synthesized according to typical procedures H and E from bicyclononene AJ2 and cyclopropyl(2,3-dimethylbenzyl)amine. LC-MS: R_t = 0.92; ES+: 632.35.

Example 623

20 **(*rac.*)-(1*R*^{*}, 5*S*^{*})-3-Acetyl-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid cyclopropyl-(3-fluoro-2-methylbenzyl)amide formate salt**

25 Synthesized according to typical procedures H and E from bicyclononene AJ2 and cyclopropyl(2-methyl-3-fluorobenzyl)amine. LC-MS: R_t = 0.91; ES+: 636.31.

Example 624

30 **(*rac.*)-(1*R*^{*}, 5*S*^{*})-3-Acetyl-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid cyclopropyl-(3,5-difluorobenzyl)amide formate salt**

Synthesized according to typical procedures H and E from bicyclononene AJ2 and cyclopropyl(3,5-difluorobenzyl)amine. LC-MS: R_t = 0.93; ES+: 672.23.

Example 625

5

(rac.)-(1R, 5S*)-3-Acetyl-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid (2-chloro-3,6-difluorobenzyl)-cyclopropylamide formate salt*

- 10 Synthesized according to typical procedures H and E from bicyclononene AJ2 and cyclopropyl(2-chloro-3,6-difluorobenzyl)amine. LC-MS: R_t = 0.90; ES+: 674.25.

Example 626

15

(rac.)-(1R, 5S*)-3-Acetyl-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid cyclopropyl-(2,3-dichlorobenzyl)amide formate salt*

- 20 Synthesized according to typical procedures H and E from bicyclononene AJ2 and cyclopropyl(2,3-dichlorobenzyl)amine. LC-MS: R_t = 0.93; ES+: 672.25.

Example 627

- 25 *(rac.)-(1R*, 5S*)-3-Acetyl-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid cyclopropyl-(3-trifluoromethoxybenzyl)amide formate salt*

- 30 Synthesized according to typical procedures H and E from bicyclononene AJ2 and cyclopropyl(3-trifluoromethoxybenzyl)amine. LC-MS: R_t = 0.93; ES+: 688.28.

Example 628

(*rac.*)-(1*R*^{*, 5*S*^{*})-3-Acetyl-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid cyclopropyl-(3-methylbenzyl)-5 amide formate salt}

Synthesized according to typical procedures H and E from bicyclononene AJ2 and cyclopropyl(3-methylbenzyl)amine. LC-MS: R_t = 0.90; ES+: 618.36.

10 Example 629

(*rac.*)-(1*R*^{*, 5*S*^{*})-3-Acetyl-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid cyclopropyl-(2,3-difluorobenzyl)amide formate salt}

15

Synthesized according to typical procedures H and E from bicyclononene AJ2 and cyclopropyl(2,3-difluorobenzyl)amine. LC-MS: R_t = 0.90; ES+: 640.29.

Example 630

20

(*rac.*)-(1*R*^{*, 5*S*^{*})-3-Acetyl-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid (3-chlorobenzyl)cyclopropyl-amide formate salt}

25

Synthesized according to typical procedures H and E from bicyclononene AJ2 and cyclopropyl(3-chlorobenzyl)amine. LC-MS: R_t = 0.91; ES+: 638.27.

Example 631

30

(*rac.*)-(1*R*^{*, 5*S*^{*})-3-Acetyl-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid cyclopropyl-(4-fluorobenzyl)-amide formate salt}

Synthesized according to typical procedures H and E from bicyclononene AJ2 and cyclopropyl(4-fluorobenzyl)amine. LC-MS: R_t = 0.89; ES+: 622.34.

5 **Example 632**

(*rac.*)-(1*R*^{*,} 5*S*^{*})-3-Acetyl-7-{4-[3-(2-bromo-5-fluorophenoxy)ethyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid (2-chlorobenzyl)-cyclopropylamide

10

Synthesized according to typical procedures H and E from bicyclononene AJ1 and (2-chlorobenzyl)cyclopropylamine. The title compound was purified by FC. LC-MS: R_t = 4.17; ES+: 666.07.

15 **Example 633**

(*rac.*)-(1*R*^{*,} 5*S*^{*})-3-Acetyl-7-{4-[3-(2-chlorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid (2-chlorobenzyl)cyclopropylamide formate salt

20

Synthesized according to typical procedures H and E from bicyclononene T2 and (2-chlorobenzyl)cyclopropylamine. LC-MS: R_t = 0.92; ES+: 617.94.

Example 634

25

(*rac.*)-(1*R*^{*,} 5*S*^{*})-3-Acetyl-7-{4-[3-(2-chlorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid [2-(4-methoxyphenoxy)ethyl]-methylamide formate salt

30 Synthesized according to typical procedures H and E from bicyclononene T2 and cyclopropyl-[2-(4-methoxyphenoxy)ethyl]amine. LC-MS: R_t = 0.84; ES+: 618.03.

Example 635

(*rac.*)-(1*R*^{*}, 5*S*^{*})-3-Acetyl-7-{4-[3-(2-chlorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid methyl-(3-trifluoromethylbenzyl)amide formate salt

Synthesized according to typical procedures H and E from bicyclononene T2 and cyclopropyl-(3-trifluoromethylbenzyl)amine. LC-MS: R_t = 0.89; ES+: 626.06.

10

Example 636

(*rac.*)-(1*R*^{*}, 5*S*^{*})-3-Acetyl-7-{4-[3-(2-chlorophenoxy)propyl]phenyl}-3,9-diaza-bicyclo[3.3.1]non-6-ene-6-carboxylic acid [2-(3,4-dimethylphenoxy)ethyl]methylamide formate salt

Synthesized according to typical procedures H and E from bicyclononene T2 and cyclopropyl-[2-(3,4-dimethylphenoxy)ethyl]amine. LC-MS: R_t = 0.91; ES+: 616.13.

20

Example 637

(*rac.*)-(1*R*^{*}, 5*S*^{*})-3-Acetyl-7-{4-[3-(2-chlorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid (3,5-dimethoxybenzyl)methylamide formate salt

Synthesized according to typical procedures H and E from bicyclononene T2 and cyclopropyl-(3,5-dimethoxybenzyl)amine. LC-MS: R_t = 0.84; ES+: 618.11.

30 **Example 638**

(*rac.*)-(1*R*^{*, 5*S*^{*})-3-Acetyl-7-[4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl]-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid (2-chlorobenzyl)cyclopropylamide trifluoroacetate salt}

5 Synthesized according to typical procedures H and E from bicyclononene AJ2 and (2-chlorobenzyl)cyclopropylamine. LC-MS: R_t = 1.00; ES+: 638.14.

Example 639

10 (*rac.*)-(1*R*^{*, 5*S*^{*})-3-Acetyl-7-[4-[3-(3-fluorophenoxy)propyl]phenyl]-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid (2-chlorobenzyl)cyclopropylamide trifluoroacetate salt}

15 Synthesized according to typical procedures H and E from bicyclononene AJ3 and (2-chlorobenzyl)cyclopropylamine. LC-MS: R_t = 0.99; ES+: 638.14.

Example 640

20 (*rac.*)-(1*R*^{*, 5*S*^{*})-5-[7-{4-[2-(2-Bromo-5-fluorophenoxy)ethyl]phenyl}-6-(methylphenethylcarbamoyl)-3,9-diazabicyclo[3.3.1]non-6-en-3-yl]-2,2-dimethyl-5-oxo-pentanoic acid trifluoroacetate salt}

25 Synthesized according to typical procedures K and E from bicyclononene AL1 and 3,3-dimethyldihydropyran-2,6-dione. LC-MS: R_t = 0.95; ES+: 710.13.

Example 641

30 (*rac.*)-(1*R*^{*, 5*S*^{*})-3-Acetyl-7-[4-[3-(2-chlorophenoxy)propyl]phenyl]-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid (2-chlorobenzyl)isopropylamide trifluoroacetate salt}

Synthesized according to typical procedures H and E from bicyclononene T2 and (2-chlorobenzyl)isopropylamine. LC-MS: R_t = 1.03; ES+: 620.20.

Example 642

5

(*rac.*)-(IR*, 5S*)-3-Acetyl-7-{4-[3-(2,3,6-trichlorophenoxy)ethyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid methylphenethylamide trifluoroacetate salt

10 Synthesized according to typical procedures F and E from bicyclononene U2 and 2,3,6-trichlorophenol. LC-MS: R_t = 1.00; ES+: 625.99.

Example 643

15 (*rac.*)-(IR*, 5S*)-3-Acetyl-7-{4-[3-(2-chlorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid (2-chlorobenzyl)ethylamide formate salt

20 Synthesized according to typical procedures H and E from bicyclononene T2 and (2-chlorobenzyl)ethylamine. LC-MS: R_t = 0.89; ES+: 606.07.

Example 644

25 (*rac.*)-(IR*, 5S*)-3-Acetyl-7-{4-[2-(3,4-dichlorophenoxy)ethoxy]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid methylphenethylamide formate salt

Synthesized according to typical procedures F and E from bicyclononene U4 and 3,4-dichlorophenol. LC-MS: R_t = 0.86; ES+: 608.01.

30

Example 645

(rac.)-(1R, 5S*)-3-Acetyl-7-{4-[2-(2-chloro-4,5-dimethylphenoxy)ethoxy]-phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid methylphenethylamide formate salt*

- 5 Synthesized according to typical procedures F and E from bicyclononene U4 and 2-chloro-4,5-dimethylphenol. LC-MS: R_t = 0.87; ES+: 602.07.

Example 646

- 10 *(rac.)-(1R*, 5S*)-3-Acetyl-7-{4-[3-(2-chlorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid (2-chlorobenzyl)methylamide formate salt*

15 Synthesized according to typical procedures F and E from bicyclononene T2 and (2-chlorobenzyl)methylamine. LC-MS: R_t = 0.87; ES+: 591.99.

Example 647

- 20 *(rac.)-(1R*, 5S*)-3-Methyl-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid (2-chlorobenzyl)cyclopropylamide formate salt*

25 A mixture of bicyclononene AL2 (1 eq.), DIPEA (3 eq.) and methyl iodide (10 eq.) in CH₂Cl₂ was stirred at rt overnight. The solvents were removed under reduced pressure and the residue treated according to typical procedure E. LC-MS: R_t = 0.89; ES+: 610.32.

Example 648

- 30 *(rac.)-(1R*, 5S*)-3-Ethyl-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid (2-chlorobenzyl)cyclopropylamide formate salt*

A mixture of bicyclononene **AL2** (1 eq.), DIPEA (3 eq.) and ethyl iodide (10 eq.) in CH₂Cl₂ was stirred at rt overnight. The solvents were removed under reduced pressure and the residue treated according to typical procedure E. LC-MS: R_t = 5 0.91; ES+: 624.33.

Example 649

(*rac.*)-(1*R*^{*}, 5*S*^{*})-3-(2-Aminoacetyl)-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]-10 phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid (2-chlorobenzyl)-cyclopropylamide formate salt

Synthesized according to typical procedures G, L and E from bicyclononene **AL2** and Boc-glycine. LC-MS: R_t = 0.80; ES+: 653.32.

15

Example 650

(*rac.*)-(1*R*^{*}, 5*S*^{*})-3-(3-Aminopropionyl)-7-{4-[3-(2,3,6-trifluorophenoxy)-propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid (2-chlorobenzyl)cyclopropylamide formate salt

Synthesized according to typical procedures G, L and E from bicyclononene **AL2** and Boc-β-alanine. LC-MS: R_t = 0.80; ES+: 667.32.

25 Example 651

(*rac.*)-(1*R*^{*}, 5*S*^{*})-3-(5-Morpholin-4-yl-5-oxopentanoyl)-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid (2-chlorobenzyl)cyclopropylamide formate salt

30

Synthesized according to typical procedure H, from bicyclononene **AL2** and glutaric anhydride, then according to typical procedure K and E from morpholine. LC-MS: R_t = 0.89; ES+: 779.31.

5 **Example 652**

(rac.)-(1*R*^{*,} 5*S*^{*})-3-(2-Tetrazol-1-ylacetyl)-7-{4-[3-(2,3,6-trifluorophenoxy)-propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid (2-chlorobenzyl)cyclopropylamide formate salt

10

Synthesized according to typical procedures G and E from bicyclononene **AL2** and (5*H*-tetrazol-5-yl)acetic acid. LC-MS: R_t = 0.91; ES+: 706.23.

Example 653

15

(rac.)-(1*R*^{*,} 5*S*^{*})-3-(5-Oxo-5-piperazin-1-ylpentanoyl)-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid (2-chlorobenzyl)cyclopropylamide formate salt

20

Synthesized according to typical procedure H, from bicyclononene **AL2** and glutaric anhydride, then according to typical procedure K, L and E from Boc-piperazine. LC-MS: R_t = 0.78; ES+: 778.37.

Example 654

25

1:1-Mixture of (1*R*, 5*S*)-3-((2*S*)-2-amino-3-hydroxypropionyl)-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid (2-chlorobenzyl)cyclopropylamide formate salt and (1*S*, 5*R*)-3-

30

((2*S*)-2-amino-3-hydroxypropionyl)-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]-phenyl}-3,9-diaza-bicyclo[3.3.1]non-6-ene-6-carboxylic acid (2-chlorobenzyl)-cyclopropylamide formate salt

Synthesized according to typical procedures G, L and E from bicyclononene **AL2** and Boc-serine. LC-MS: R_t = 0.79; ES+: 683.34.

Example 655

5

1:1-Mixture of (*1R*, *5S*)-3-((2*S*)-2-aminopropionyl)-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid (2-chlorobenzyl)cyclopropylamide formate salt and (*1S*, *5R*)-3-((2*S*)-2-aminopropionyl)-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid (2-chlorobenzyl)cyclopropylamide formate salt

10

Synthesized according to typical procedures G, L and E from bicyclononene **AL2** and Boc-alanine. LC-MS: R_t = 0.80; ES+: 667.32.

15

Example 656

(*rac.*)-(*1R*^{*}, *5S*^{*})-3-Acetyl-7-{4-[2-(2,6-dichloro-4-methylphenoxy)ethoxy]-phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid cyclopropyl-(2,3-dichlorobenzyl)amide formate salt

20

Synthesized according to typical procedures H and E from bicyclononene **BN4** and (2,3-dichlorobenzyl)cyclopropylamine. LC-MS: R_t = 0.96; ES+: 702.09.

25

Example 657

(*rac.*)-(*1R*^{*}, *5S*^{*})-3-Acetyl-7-{4-[2-(2,6-dichloro-4-methylphenoxy)ethoxy]-phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid cyclopropyl-(2,3-dimethylbenzyl)amide formate salt

30

Synthesized according to typical procedures H and E from bicyclononene **BN4** and cyclopropyl-(2,3-dimethylbenzyl)amine. LC-MS: R_t = 0.94; ES+: 662.27.

Example 658

(*rac.*)-(1*R*^{*}, 5*S*^{*})-3-Acetyl-7-{4-[2-(2,6-dichloro-4-methylphenoxy)ethoxy]-phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid (2-chloro-3-trifluoromethylbenzyl)cyclopropylamide formate salt

Synthesized according to typical procedures H and E from bicyclononene BN4 and (2-chloro-3-trifluoromethylbenzyl)cyclopropylamine. LC-MS: R_t = 0.96; ES+: 736.08.

Example 659

(*rac.*)-(1*R*^{*}, 5*S*^{*})-3-Acetyl-7-{4-[2-(2,6-dichloro-4-methylphenoxy)ethoxy]-phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid (2-bromobenzyl)-cyclopropylamide formate salt

Synthesized according to typical procedures H and E from bicyclononene BN4 and (2-bromobenzyl)cyclopropylamine. LC-MS: R_t = 0.92; ES+: 682.20.

20

Example 660

(*rac.*)-(1*R*^{*}, 5*S*^{*})-3-Acetyl-7-{4-[2-(2,6-dichloro-4-methylphenoxy)ethoxy]-phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid (2-chlorobenzyl)-cyclopropylamide formate salt

Synthesized according to typical procedures H and E from bicyclononene BN4 and (2-chlorobenzyl)cyclopropylamine. LC-MS: R_t = 0.93; ES+: 668.20.

30 **Example 661**

(*rac.*)-(1*R*^{*}, 5*S*^{*})-3-Acetyl-7-{4-[2-(2,6-dichloro-4-methylphenoxy)ethoxy]-phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid (2-chlorobenzyl)-ethylamide formate salt

- 5 Synthesized according to typical procedures H and E from bicyclononene BN4 and (2-chlorobenzyl)ethylamine. LC-MS: R_t = 0.93; ES+: 656.20.

Example 662

10 **(*rac.*)-(1*R*^{*}, 5*S*^{*})-3-Acetyl-7-{4-[2-(2,6-dichloro-4-methylphenoxy)ethoxy]-phenyl}-3,9-diaza-bicyclo[3.3.1]non-6-ene-6-carboxylic acid (6-chlorobenzo-[1,3]dioxol-5-ylmethyl)cyclopropylamide formate salt**

- 15 Synthesized according to typical procedures H and E from bicyclononene BN4 and (6-chlorobenzo[1,3]dioxol-5-ylmethyl)cyclopropylamine. LC-MS: R_t = 0.93; ES+: 712.15.

Example 663

20 **(*rac.*)-(1*R*^{*}, 5*S*^{*})-3-Acetyl-7-{4-[2-(2,6-dichloro-4-methylphenoxy)ethoxy]-phenyl}-3,9-diaza-bicyclo[3.3.1]non-6-ene-6-carboxylic acid cyclopropyl-(3,5-dimethoxybenzyl)amide formate salt**

- 25 Synthesized according to typical procedures H and E from bicyclononene BN4 and cyclopropyl-(3,5-dimethoxybenzyl)amine. LC-MS: R_t = 0.92; ES+: 694.20.

Example 664

30 **(*rac.*)-(1*R*^{*}, 5*S*^{*})-3-Acetyl-7-{4-[2-(2,6-dichloro-4-methylphenoxy)ethoxy]-phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid cyclopropyl-(3-methoxybenzyl)amide formate salt**

Synthesized according to typical procedures H and E from bicyclononene BN4 and cyclopropyl-(3-methoxybenzyl)amine. LC-MS: R_t = 0.91; ES+: 664.25.

Example 665

5

(*rac.*)-(1*R*^{*, 5*S*^{*})-3-Acetyl-7-{4-[2-(2-chloro-5-fluorophenoxy)ethoxy]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid cyclopropyl-(2,3-dichlorobenzyl)amide formate salt}

10 Synthesized according to typical procedures H and E from bicyclononene BN10 and cyclopropyl-(2,3-dichlorobenzyl)amine. LC-MS: R_t = 0.92; ES+: 672.22.

Example 666

15 (*rac.*)-(1*R*^{*, 5*S*^{*})-3-Acetyl-7-{4-[2-(2-chloro-5-fluorophenoxy)ethoxy]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid (2-chloro-3-trifluoromethylbenzyl)cyclopropylamide formate salt}

20 Synthesized according to typical procedures H and E from bicyclononene BN10 and (2-chloro-3-trifluoromethylbenzyl)cyclopropylamine. LC-MS: R_t = 0.93; ES+: 706.09.

Example 667

25 (*rac.*)-(1*R*^{*, 5*S*^{*})-3-Acetyl-7-{4-[2-(2,6-dichloro-4-methylphenoxy)ethoxy]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid (3-chlorobenzyl)cyclopropylamide formate salt}

30 Synthesized according to typical procedures H and E from bicyclononene BN4 and (3-chlorobenzyl)cyclopropylamine. LC-MS: R_t = 0.93; ES+: 668.21.

Example 668

(*rac.*)-(IR*, 5S*)-3-Acetyl-7-{4-[2-(2,6-dichloro-4-methylphenoxy)ethoxy]-phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid cyclopropyl-(3-methylbenzyl)amide formate salt

5

Synthesized according to typical procedures H and E from bicyclononene BN4 and cyclopropyl-(3-methylbenzyl)amine. LC-MS: R_t = 0.93; ES+: 648.23.

Example 669

10

(*rac.*)-(IR*, 5S*)-3-Acetyl-7-{4-[2-(2,6-dichloro-4-methylphenoxy)ethoxy]-phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid cyclopropyl-(2-fluoro-5-methoxybenzyl)amide formate salt

15 Synthesized according to typical procedures H and E from bicyclononene BN4 and cyclopropyl-(2-fluoro-5-methoxybenzyl)amine. LC-MS: R_t = 0.92; ES+: 682.20.

Example 670

20

(*rac.*)-(IR*, 5S*)-3-Acetyl-7-{4-[2-(2-chloro-4,5-dimethylphenoxy)-ethoxy]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid cyclopropyl-(2,3-dichlorobenzyl)amide formate salt

25 Synthesized according to typical procedures H and E from bicyclononene BN3 and cyclopropyl-(2,3-dichlorobenzyl)amine. LC-MS: R_t = 0.95.; ES+: 684.20.

Example 671

30 **(*rac.*)-(IR*, 5S*)-3-Acetyl-7-{4-[3-(2-chloro-3,6-difluorophenoxy)propyl]-phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid (2-chlorobenzyl)-cyclopropylamide formate salt**

Synthesized according to typical procedures F and E from bicyclononene BB1 and 2-chloro-3,6-difluorophenol. LC-MS: R_t = 0.93; ES+: 654.28.

5 **Example 672**

(*rac.*)-(1*R*^{*}, 5*S*^{*})-3-Acetyl-7-{4-[2-(2-chloro-4,5-dimethylphenoxy)ethoxy]-phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid (2-chloro-3-trifluoromethylbenzyl)cyclopropylamide formate salt

10

Synthesized according to typical procedures H and E from bicyclononene BN3 and (2-chloro-3-trifluoromethylbenzyl)cyclopropylamine. LC-MS: R_t = 0.96; ES+: 716.18.

15 **Example 673**

(*rac.*)-(1*R*^{*}, 5*S*^{*})-3-Acetyl-7-{4-[2-(2,6-dichloro-4-methylphenoxy)ethoxy]-phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid cyclopropyl-(3,4-dimethoxybenzyl)amide formate salt

20

Synthesized according to typical procedures H and E from bicyclononene BN4 and cyclopropyl-(3,4-dimethoxybenzyl)amine. LC-MS: R_t = 0.89; ES+: 694.23.

Example 674

25

(*rac.*)-(1*R*^{*}, 5*S*^{*})-3-Acetyl-7-{4-[2-(2,6-dichloro-4-methylphenoxy)ethoxy]-phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid cyclopropyl-(3-trifluoromethoxybenzyl)amide formate salt

30

Synthesized according to typical procedures H and E from bicyclononene BN4 and cyclopropyl-(3-trifluoromethoxybenzyl)amine. LC-MS: R_t = 0.95; ES+: 718.13.

Example 675

(*rac.*)-(IR*, 5S*)-3-Acetyl-7-{4-[2-(2,4,5-trichlorophenoxy)ethoxy]phenyl}-
5 3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid (2-chloro-3-trifluoro-
methylbenzyl)cyclopropylamide formate salt

Synthesized according to typical procedures H and E from bicyclononene BN9
and (2-chloro-3-trifluoromethylbenzyl)cyclopropylamine. LC-MS: R_t = 0.97;
10 ES+: 758.08.

Example 676

(*rac.*)-(IR*, 5S*)-3-Acetyl-7-{4-[2-(2,4,5-trichlorophenoxy)ethoxy]phenyl}-
15 3,9-diazabicyclo-[3.3.1]non-6-ene-6-carboxylic acid cyclopropyl-(2,3-
dichlorobenzyl)amide formate salt

Synthesized according to typical procedures H and E from bicyclononene BN9
and cyclopropyl-(2,3-dichlorobenzyl)amine. LC-MS: R_t = 0.96; ES+: 724.02.
20

Example 677

(*rac.*)-(IR*, 5S*)-3-Acetyl-7-{4-[2-(2,3-dichlorophenoxy)ethoxy]phenyl}-3,9-
diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid cyclopropyl-(2,3-dimethyl-
25 benzyl)amide formate salt

Synthesized according to typical procedures H and E from bicyclononene BN5
and cyclopropyl-(2,3-dimethylbenzyl)amine. LC-MS: R_t = 0.92; ES+: 648.25.

30 Example 678

(*rac.*)-(1*R*^{*}, 5*S*^{*})-3-Acetyl-7-{4-[2-(2,4,5-trichlorophenoxy)ethoxy]phenyl}-3,9-diazabicyclo-[3.3.1]non-6-ene-6-carboxylic acid cyclopropyl-(2,3-dimethylbenzyl)amide formate salt

5 Synthesized according to typical procedures H and E from bicyclononene BN9 and cyclopropyl(2,3-dimethylbenzyl)amine. LC-MS: R_t = 0.95; ES+: 684.12.

Example 679

10 **(*rac.*)-(1*R*^{*}, 5*S*^{*})-3-Acetyl-7-{4-[2-(2-chloro-4,5-dimethylphenoxy)ethoxy]-phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid cyclopropyl-(2,3-dimethylbenzyl)amide formate salt**

15 Synthesized according to typical procedures H and E from bicyclononene BN3 and cyclopropyl-(2,3-dimethylbenzyl)amine. LC-MS: R_t = 0.94; ES+: 642.31.

Example 680

20 **(*rac.*)-(1*R*^{*}, 5*S*^{*})-3-Acetyl-7-{4-[2-(2-bromo-5-fluorophenoxy)ethoxy]phenyl}-3,9-diaza-bicyclo[3.3.1]non-6-ene-6-carboxylic acid cyclopropyl-(2,3-dichlorobenzyl)amide formate salt**

25 Synthesized according to typical procedures H and E from bicyclononene BN2 and cyclopropyl-(2,3-dichlorobenzyl)amine. LC-MS: R_t = 0.92; ES+: 718.05.

25

Example 681

30 **(*rac.*)-(1*R*^{*}, 5*S*^{*})-3-Acetyl-7-{4-[2-(2-chloro-4,5-dimethylphenoxy)ethoxy]-phenyl}-3,9-diaza-bicyclo[3.3.1]non-6-ene-6-carboxylic acid (2-bromobenzyl)-cyclopropylamide formate salt**

Synthesized according to typical procedures H and E from bicyclononene BN3 and (2-bromobenzyl)cyclopropylamine. LC-MS: R_t = 0.94; ES+: 694.15.

Example 682

5

(*rac.*)-(1*R*^{*,}, 5*S*^{*})-3-Acetyl-7-{4-[2-(2,3-dichlorophenoxy)ethoxy]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid cyclopropyl-(2,3-dichlorobenzyl)amide formate salt

10 Synthesized according to typical procedures H and E from bicyclononene BN5 and cyclopropyl-(2,3-dichlorobenzyl)amine. LC-MS: R_t = 0.93; ES+: 690.06.

Example 683

15 (*rac.*)-(1*R*^{*,}, 5*S*^{*})-3-Acetyl-7-{4-[2-(2-chloro-5-fluorophenoxy)ethoxy]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid cyclopropyl-(2,3-dimethylbenzyl)amide formate salt

20 Synthesized according to typical procedures H and E from bicyclononene BN10 and cyclopropyl-(2,3-dimethylbenzyl)amine. LC-MS: R_t = 0.91; ES+: 632.31.

Example 684

25 (*rac.*)-(1*R*^{*,}, 5*S*^{*})-3-Acetyl-7-{4-[2-(2,4,5-trichlorophenoxy)ethoxy]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid (2-chlorobenzyl)-cyclopropylamide formate salt

Synthesized according to typical procedures H and E from bicyclononene BN9 and (2-chlorobenzyl)cyclopropylamine. LC-MS: R_t = 0.94; ES+: 690.07.

30

Example 685

(*rac.*)-(1*R*^{*}, 5*S*^{*})-3-Acetyl-7-{4-[2-(2-bromo-5-fluorophenoxy)ethoxy]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid (2-chloro-3-trifluoromethylbenzyl)cyclopropylamide formate salt

- 5 Synthesized according to typical procedures H and E from bicyclononene BN2 and (2-chloro-3-trifluoromethylbenzyl)cyclopropylamine. LC-MS: R_t = 0.93; ES+: 752.06.

Example 686

10

(*rac.*)-(1*R*^{*}, 5*S*^{*})-3-Acetyl-7-{4-[2-(2,4,5-trichlorophenoxy)ethoxy]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid (2-bromobenzyl)cyclopropylamide formate salt

- 15 Synthesized according to typical procedures H and E from bicyclononene BN9 and (2-bromobenzyl)cyclopropylamine. LC-MS: R_t = 0.95; ES+: 733.99.

Example 687

- 20 **(*rac.*)-(1*R*^{*}, 5*S*^{*})-3-Acetyl-7-{4-[2-(2-bromo-5-fluorophenoxy)ethoxy]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid cyclopropyl-(2,3-dimethylbenzyl)amide formate salt**

25 Synthesized according to typical procedures H and E from bicyclononene BN2 and cyclopropyl-(2,3-dimethylbenzyl)amine. LC-MS: R_t = 0.91; ES+: 678.22.

Example 688

- 30 **(*rac.*)-(1*R*^{*}, 5*S*^{*})-3-Acetyl-7-{4-[2-(2,3-dichlorophenoxy)ethoxy]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid (2-chloro-3-trifluoromethylbenzyl)cyclopropylamide formate salt**

Synthesized according to typical procedures H and E from bicyclononene BN5 and (2-chloro-3-trifluoromethylbenzyl)cyclopropylamine. LC-MS: R_t = 0.94; ES+: 724.13.

5 **Example 689**

(*rac.*)-(1*R*^{*,}, 5*S*^{*)}-3-Acetyl-7-{4-[2-(2-chloro-4,5-dimethylphenoxy)ethoxy]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid (3-chlorobenzyl)cyclopropylamide formate salt

10

Synthesized according to typical procedures H and E from bicyclononene BN3 and (3-chlorobenzyl)cyclopropylamine. LC-MS: R_t = 0.93; ES+: 648.26.

Example 690

15

(*rac.*)-(1*R*^{*,}, 5*S*^{*)}-3-Acetyl-7-{4-[2-(2,4,5-trichlorophenoxy)ethoxy]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid (2-chlorobenzyl)ethylamide formate salt

20

Synthesized according to typical procedures H and E from bicyclononene BN9 and (2-chlorobenzyl)ethylamine. LC-MS: R_t = 0.94; ES+: 678.12.

Example 691

25

(*rac.*)-(1*R*^{*,}, 5*S*^{*)}-3-Acetyl-7-{4-[2-(2-chloro-5-fluorophenoxy)ethoxy]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid (2-bromobenzyl)cyclopropylamide formate salt

30

Synthesized according to typical procedures H and E from bicyclononene BN10 and (2-bromobenzyl)cyclopropylamine. LC-MS: R_t = 0.90; ES+: 684.11.

Example 692

(rac.)-(IR, 5S*)-3-Acetyl-7-{4-[2-(4-chloro-2-methylphenoxy)ethoxy]-phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid cyclopropyl-(2,3-dichlorobenzyl)amide formate salt*

5

Synthesized according to typical procedures H and E from bicyclononene BN7 and cyclopropyl-(2,3-dichlorobenzyl)amine. LC-MS: R_t = 0.94; ES+: 670.17

Example 693

10

(rac.)-(IR, 5S*)-3-Acetyl-7-{4-[2-(2-chloro-4,5-dimethylphenoxy)ethoxy]-phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid cyclopropyl-(3-methoxybenzyl)amide formate salt*

15 Synthesized according to typical procedures H and E from bicyclononene BN3 and cyclopropyl-(3-methoxybenzyl)amine. LC-MS: R_t = 0.91; ES+: 644.32.

Example 694

20 *(rac.)-(IR*, 5S*)-3-Acetyl-7-{4-[2-(2-bromo-5-fluorophenoxy)ethoxy]-phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid (2-bromobenzyl)-cyclopropylamide formate salt*

25 Synthesized according to typical procedures H and E from bicyclononene BN2 and cyclopropyl(2-bromobenzyl)amine. LC-MS: R_t = 0.91; ES+: 728.04.

Example 695

30 *(rac.)-(IR*, 5S*)-3-Acetyl-7-{4-[2-(4-chloro-2-methylphenoxy)ethoxy]-phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid cyclopropyl-(2,3-dimethylbenzyl)amide formate salt*

Synthesized according to typical procedures H and E from bicyclononene BN7 and cyclopropyl(2,3-dimethylbenzyl)amine. LC-MS: R_t = 0.93; ES+: 628.30.

Example 696

5

(*rac.*)-(1*R*^{*,}, 5*S*^{*})-3-Acetyl-7-{4-[2-(2,4,5-trichlorophenoxy)ethoxy]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid cyclopropyl-(3-methoxybenzyl)amide formate salt

10 Synthesized according to typical procedures H and E from bicyclononene BN9 and cyclopropyl-(3-methoxybenzyl)amine. LC-MS: R_t = 0.92; ES+: 686.14.

Example 697

15 (*rac.*)-(1*R*^{*,}, 5*S*^{*})-3-Acetyl-7-{4-[2-(2-chloro-4,5-dimethylphenoxy)ethoxy]-phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid cyclopropyl-(2-fluoro-5-methoxybenzyl)amide formate salt

20 Synthesized according to typical procedures H and E from bicyclononene BN3 and cyclopropyl-(2-fluoro-5-methoxybenzyl)amine. LC-MS: R_t = 0.91; ES+: 662.29.

Example 698

25 (*rac.*)-(1*R*^{*,}, 5*S*^{*})-3-Acetyl-7-{4-[2-(4-chloro-2-methylphenoxy)ethoxy]-phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid (2-bromobenzyl)-cyclopropylamide formate salt

30 Synthesized according to typical procedures H and E from bicyclononene BN7 and (2-bromobenzyl)cyclopropylamine. LC-MS: R_t = 0.93; ES+: 680.15.

Example 699

(*rac.*)-(1*R*^{*, 5*S*})-3-Acetyl-7-{4-[2-(2-chloro-5-fluorophenoxy)ethoxy]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid (2-chlorobenzyl)cyclopropylamide formate salt

5

Synthesized according to typical procedures H and E from bicyclononene BN10 and (2-chlorobenzyl)cyclopropylamine. LC-MS: R_t = 0.90; ES+: 638.22.

Example 700

10

(*rac.*)-(1*R*^{*, 5*S*})-3-Acetyl-7-{4-[2-(2,4,5-trichlorophenoxy)ethoxy]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid cyclopropyl-(3,5-dimethoxybenzyl)amide formate salt

15

Synthesized according to typical procedures H and E from bicyclononene BN9 and cyclopropyl-(3,5-dimethoxybenzyl)amine. LC-MS: R_t = 0.93; ES+: 716.14.

Example 701

20

(*rac.*)-(1*R*^{*, 5*S*})-3-Acetyl-7-{4-[2-(2,3-dichlorophenoxy)ethoxy]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid (2-bromobenzyl)cyclopropylamide formate salt

25

Synthesized according to typical procedures H and E from bicyclononene BN5 and (2-bromobenzyl)cyclopropylamine. LC-MS: R_t = 0.91; ES+: 700.07.

Example 702

30

(*rac.*)-(1*R*^{*, 5*S*})-3-Acetyl-7-{4-[2-(2-chloro-4,5-dimethylphenoxy)ethoxy]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid (6-chlorobenzo[1,3]dioxol-5-ylmethyl)-cyclopropylamide formate salt

Synthesized according to typical procedures H and E from bicyclononene BN3 and (6-chlorobenzo[1,3]dioxol-5-ylmethyl)cyclopropylamine. LC-MS: R_t = 0.93; ES+: 694.17.

5 **Example 703**

(*rac.*)-(IR*, 5S*)-3-Acetyl-7-{4-[2-(2,3-dichlorophenoxy)ethoxy]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid (2-chlorobenzyl)cyclopropylamide formate salt

10

Synthesized according to typical procedures H and E from bicyclononene BN5 and (2-chlorobenzyl)cyclopropylamine. LC-MS: R_t = 0.91; ES+: 656.19.

Example 704

15

(*rac.*)-(IR*, 5S*)-3-Acetyl-7-{4-[2-(4-chloro-2-methylphenoxy)ethoxy]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid (2-chloro-3-trifluoromethylbenzyl)cyclopropylamide formate salt

20

Synthesized according to typical procedures H and E from bicyclononene BN7 and (2-chloro-3-trifluoromethylbenzyl)cyclopropylamine. LC-MS: R_t = 0.95; ES+: 702.17.

Example 705

25

(*rac.*)-(IR*, 5S*)-3-Acetyl-7-{4-[2-(2-chloro-4,5-dimethylphenoxy)ethoxy]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid cyclopropyl-(3-methylbenzyl)amide formate salt

30

Synthesized according to typical procedures H and E from bicyclononene BN3 and cyclopropyl-(3-methylbenzyl)amine. LC-MS: R_t = 0.92; ES+: 628.32.

Example 706

(*rac.*)-(IR*, 5S*)-3-Acetyl-7-{4-[2-(2,3-dichlorophenoxy)ethoxy]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid (2-chlorobenzyl)ethylamide
5 formate salt

Synthesized according to typical procedures H and E from bicyclononene BN5 and (2-chlorobenzyl)ethylamine. LC-MS: R_t = 0.90; ES+: 642.19.

10 Example 707

(*rac.*)-(IR*, 5S*)-3-Acetyl-7-{4-[2-(2-chloro-5-fluorophenoxy)ethoxy]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid cyclopropyl-(2-fluoro-5-methoxybenzyl)amide formate salt

15

Synthesized according to typical procedures H and E from bicyclononene BN10 and cyclopropyl-(2-fluoro-5-methoxybenzyl)amine. LC-MS: R_t = 0.88; ES+: 652.26.

20 Example 708

(*rac.*)-(IR*, 5S*)-3-Acetyl-7-{4-[2-(4-chloro-2-methylphenoxy)ethoxy]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid (2-chlorobenzyl)-cyclopropylamide formate salt

25

Synthesized according to typical procedures H and E from bicyclononene BN7 and cyclopropyl(2-chlorobenzyl)amine. LC-MS: R_t = 0.92; ES+: 634.22.

Example 709

30

(rac.)-(1R, 5S*)-3-Acetyl-7-{4-[2-(2-bromo-5-fluorophenoxy)ethoxy]-phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid cyclopropyl-(3,5-dimethoxybenzyl)amide formate salt*

5 Synthesized according to typical procedures H and E from bicyclononene BN2 and cyclopropyl-(3,5-dimethoxybenzyl)amine. LC-MS: R_t = 0.88; ES+: 708.14.

Example 710

10 *(rac.)-(1R*, 5S*)-3-Acetyl-7-{4-[2-(2-chloro-5-fluorophenoxy)ethoxy]-phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid cyclopropyl-(3-methoxybenzyl)amide formate salt*

15 Synthesized according to typical procedures H and E from bicyclononene BN10 and cyclopropyl-(3-methoxybenzyl)amine. LC-MS: R_t = 0.87; ES+: 634.27.

Example 711

20 *(rac.)-(1R*, 5S*)-3-Acetyl-7-{4-[2-(2,4,5-trichlorophenoxy)ethoxy]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid (3-chlorobenzyl)-cyclopropylamide formate salt*

25 Synthesized according to typical procedures H and E from bicyclononene BN9 and (3-chlorobenzyl)cyclopropylamine. LC-MS: R_t = 0.94; ES+: 690.09.

Example 712

30 *(rac.)-(1R*, 5S*)-3-Acetyl-7-{4-[2-(2,6-dichloro-4-methylphenoxy-ethoxy]-phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid cyclopropyl-(3,5-difluorobenzyl)amide formate salt*

Synthesized according to typical procedures H and E from bicyclononene **BN4** and cyclopropyl-(3,5-difluorobenzyl)amine. LC-MS: R_t = 0.93; ES+: 670.22.

Example 713

5

(*rac.*)-(1*R*^{*}, 5*S*^{*})-3-Acetyl-7-{4-[2-(2,4,5-trichlorophenoxy)ethoxy]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid cyclopropyl-(2-fluoro-5-methoxybenzyl)amide formate salt

- 10 Synthesized according to typical procedures H and E from bicyclononene **BN9** and cyclopropyl-(2-fluoro-5-methoxybenzyl)amine. LC-MS: R_t = 0.93; ES+: 702.40.

Example 714

15

(*rac.*)-(1*R*^{*}, 5*S*^{*})-3-Acetyl-7-{4-[2-(2,4,5-trichlorophenoxy)ethoxy]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid cyclopropyl-(3,4-dimethoxybenzyl)amide formate salt

- 20 Synthesized according to typical procedures H and E from bicyclononene **BN9** and cyclopropyl-(3,4-dimethoxybenzyl)amine. LC-MS: R_t = 0.90; ES+: 716.10.

Example 715

- 25 (*rac.*)-(1*R*^{*}, 5*S*^{*})-3-Acetyl-7-{4-[2-(2-chloro-5-fluorophenoxy)ethoxy]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid (6-chloro-benzo[1,3]-dioxol-5-ylmethyl)cyclopropylamide formate salt

- 30 Synthesized according to typical procedures H and E from bicyclononene **BN10** and (6-chlorobenzo[1,3]dioxol-5-ylmethyl)cyclopropylamine. LC-MS: R_t = 0.90; ES+: 684.19.

Example 716

(*rac.*)-(IR*, 5S*)-3-Acetyl-7-{4-[2-(2-bromo-5-fluorophenoxy)ethoxy]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid (3-chlorobenzyl)-5-cyclopropylamide formate salt

Synthesized according to typical procedures H and E from bicyclononene BN2 and (3-chlorobenzyl)cyclopropylamine. LC-MS: R_t = 0.90; ES+: 682.16.

10 Example 717

(*rac.*)-(IR*, 5S*)-3-Acetyl-7-{4-[2-(2,3-dichlorophenoxy)ethoxy]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid cyclopropyl-(3,4-dimethoxybenzyl)amide formate salt

15

Synthesized according to typical procedures H and E from bicyclononene BN5 and cyclopropyl-(3,4-dimethoxybenzyl)amine. LC-MS: R_t = 0.97; ES+: 758.08.

Example 718

20

(*rac.*)-(IR*, 5S*)-3-Acetyl-7-{4-[2-(2-chloro-5-fluorophenoxy)ethoxy]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid (2-chlorobenzyl)-ethylamide formate salt

25

Synthesized according to typical procedures H and E from bicyclononene BN10 and (2-chlorobenzyl)ethylamine. LC-MS: R_t = 0.89; ES+: 626.25.

Example 719

30

(*rac.*)-(IR*, 5S*)-3-Acetyl-7-{4-[2-(4-chloro-2-methylphenoxy)ethoxy]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid cyclopropyl-(3,5-dimethoxybenzyl)amide formate salt

Synthesized according to typical procedures H and E from bicyclononene BN7 and cyclopropyl-(3,5-dimethoxybenzyl)amine. LC-MS: R_t = 0.90; ES+: 660.29.

5 **Example 720**

(rac.)-(IR*, 5S*)-3-Acetyl-7-{4-[2-(2,4,5-trichlorophenoxy)ethoxy]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid cyclopropyl-(3-methylbenzyl)amide formate salt

10

Synthesized according to typical procedures H and E from bicyclononene BN9 and cyclopropyl-(3-methylbenzyl)amine. LC-MS: R_t = 0.94; ES+: 670.22.

15 **Example 721**

(rac.)-(IR*, 5S*)-3-Acetyl-7-{4-[2-(2,3-dichlorophenoxy)ethoxy]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid cyclopropyl-(2-fluoro-5-methoxybenzyl)amide formate salt

20 Synthesized according to typical procedures H and E from bicyclononene BN5 and cyclopropyl-(2-fluoro-5-methoxybenzyl)amine. LC-MS: R_t = 0.89; ES+: 668.25.

25 **Example 722**

(rac.)-(IR*, 5S*)-3-Acetyl-7-{4-[2-(2-chloro-4,5-dimethylphenoxy)ethoxy]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid cyclopropyl-(3,5-difluorobenzyl)amide formate salt

30 Synthesized according to typical procedures H and E from bicyclononene BN3 and cyclopropyl-(3,5-difluorobenzyl)amine. LC-MS: R_t = 0.92; ES+: 650.23.

Example 723

5 **(*rac.*)-(IR*, 5S*)-3-Acetyl-7-{4-[2-(2,3-dichlorophenoxy)ethoxy]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid cyclopropyl-(3,5-dimethoxybenzyl)amide formate salt**

Synthesized according to typical procedures H and E from bicyclononene BN5 and cyclopropyl-(3,5-dimethoxybenzyl)amine. LC-MS: R_t = 0.89; ES+: 680.22.

10 **Example 724**

15 **(*rac.*)-(IR*, 5S*)-3-Acetyl-7-{4-[2-(2-chloro-5-fluorophenoxy)ethoxy]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid (2-chlorobenzyl)-cyclopropylamide formate salt**

Synthesized according to typical procedures H and E from bicyclononene BN10 and cyclopropyl(2-chlorobenzyl)amine. LC-MS: R_t = 0.90; ES+: 638.22.

20 **Example 725**

25 **(*rac.*)-(IR*, 5S*)-3-Acetyl-7-{4-[2-(2-bromo-5-fluorophenoxy)ethoxy]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid (2-chlorobenzyl)-cyclopropylamide formate salt**

30 Synthesized according to typical procedures H and E from bicyclononene BN2 and (2-chlorobenzyl)cyclopropylamine. LC-MS: R_t = 0.90; ES+: 684.12.

Example 726

35 **(*rac.*)-(IR*, 5S*)-3-Acetyl-7-{4-[2-(2-chloro-4,5-dimethylphenoxy)ethoxy]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid cyclopropyl-(3,4-dimethoxybenzyl)amide formate salt**

Synthesized according to typical procedures H and E from bicyclononene BN3 and cyclopropyl-(3,4-dimethoxybenzyl)amine. LC-MS: R_t = 0.88; ES+: 674.31.

5 **Example 727**

(*rac.*)-(1*R*^{*}, 5*S*^{*})-3-Acetyl-7-{4-[2-(2-chloro-4,5-dimethylphenoxy)ethoxy]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid cyclopropyl-(3-trifluoromethoxybenzyl)amide formate salt

10

Synthesized according to typical procedures H and E from bicyclononene BN3 and cyclopropyl-(3-trifluoromethoxybenzyl)amine. LC-MS: R_t = 0.95; ES+: 698.22.

15 **Example 728**

(*rac.*)-(1*R*^{*}, 5*S*^{*})-3-Acetyl-7-{4-[2-(2-chloro-5-fluorophenoxy)ethoxy]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid cyclopropyl-(3-methylbenzyl)amide formate salt

20

Synthesized according to typical procedures H and E from bicyclononene BN10 and cyclopropyl-(3-methylbenzyl)amine. LC-MS: R_t = 0.89; ES+: 618.28.

Example 729

25

(*rac.*)-(1*R*^{*}, 5*S*^{*})-3-Acetyl-7-{4-[2-(2-chloro-5-fluorophenoxy)ethoxy]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid cyclopropyl-(3,5-dimethoxybenzyl)amide formate salt

30

Synthesized according to typical procedures H and E from bicyclononene BN10 and cyclopropyl-(3,5-dimethoxybenzyl)amine. LC-MS: R_t = 0.88; ES+: 664.27.

Example 730

(*rac.*)-(1*R*^{*,}, 5*S*^{*})-3-Acetyl-7-{4-[2-(2,3-dichlorophenoxy)ethoxy]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid cyclopropyl-(3-methylbenzyl)amide formate salt

Synthesized according to typical procedures H and E from bicyclononene BN5 and cyclopropyl-(3-methylbenzyl)amine. LC-MS: R_t = 0.90; ES+: 634.21.

10 **Example 731**

(*rac.*)-(1*R*^{*,}, 5*S*^{*})-3-Acetyl-7-{4-[2-(2-bromo-5-fluorophenoxy)ethoxy]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid cyclopropyl-(2-fluoro-5-methoxybenzyl)amide formate salt

15 Synthesized according to typical procedures H and E from bicyclononene BN2 and cyclopropyl-(2-fluoro-5-methoxybenzyl)amine. LC-MS: R_t = 0.89; ES+: 696.14.

20 **Example 732**

(*rac.*)-(1*R*^{*,}, 5*S*^{*})-3-Acetyl-7-{4-[2-(2-bromo-5-fluorophenoxy)ethoxy]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid (2-chlorobenzyl)ethylamide formate salt

25 Synthesized according to typical procedures H and E from bicyclononene BN2 and (2-chlorobenzyl)ethylamine. LC-MS: R_t = 0.90; ES+: 672.14.

Example 733

(*rac.*)-(1*R*^{*}, 5*S*^{*})-3-Acetyl-7-{4-[2-(2,3-dichlorophenoxy)ethoxy]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid (6-chlorobenzo[1,3]dioxol-5-ylmethyl)cyclopropylamide formate salt

5 Synthesized according to typical procedures H and E from bicyclononene BN5 and (6-chlorobenzo[1,3]dioxol-5-ylmethyl)cyclopropylamine. LC-MS: R_t = 0.91; ES+: 700.12.

Example 734

10

(*rac.*)-(1*R*^{*}, 5*S*^{*})-3-Acetyl-7-{4-[2-(2,3-dichlorophenoxy)ethoxy]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid (3-chlorobenzyl)cyclopropylamide formate salt

15 Synthesized according to typical procedures H and E from bicyclononene BN5 and (3-chlorobenzyl)cyclopropylamine. LC-MS: R_t = 0.91; ES+: 656.19.

Example 735

20 (*rac.*)-(1*R*^{*}, 5*S*^{*})-3-Acetyl-7-{4-[2-(4-chloro-2-methylphenoxy)ethoxy]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid (6-chlorobenzo[1,3]dioxol-5-ylmethyl)cyclopropylamide formate salt

25 Synthesized according to typical procedures H and E from bicyclononene BN7 and (6-chlorobenzo[1,3]dioxol-5-ylmethyl)cyclopropylamine. LC-MS: R_t = 0.92; ES+: 678.20.

Example 736

30 (*rac.*)-(1*R*^{*}, 5*S*^{*})-3-Acetyl-7-{4-[2-(2-bromo-5-fluorophenoxy)ethoxy]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid cyclopropyl-(3-methoxybenzyl)amide formate salt

Synthesized according to typical procedures H and E from bicyclononene BN2 and cyclopropyl-(3-methoxybenzyl)amine. LC-MS: R_t = 0.88; ES+: 678.20.

5 **Example 737**

(*rac.*)-(IR*, 5S*)-3-Acetyl-7-{4-[2-(4-chloro-2-methylphenoxy)ethoxy]-phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid (2-chlorobenzyl)-ethylamide formate salt

10

Synthesized according to typical procedures H and E from bicyclononene BN7 and (2-chlorobenzyl)ethylamine. LC-MS: R_t = 0.91; ES+: 622.26.

Example 738

15

(*rac.*)-(IR*, 5S*)-3-Acetyl-7-{4-[2-(4-chloro-2-methylphenoxy)ethoxy]-phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid cyclopropyl-(3-methylbenzyl)amide formate salt

20

Synthesized according to typical procedures H and E from bicyclononene BN7 and cyclopropyl-(3-methylbenzyl)amine. LC-MS: R_t = 0.91; ES+: 614.32.

Example 739

25

(*rac.*)-(IR*, 5S*)-3-Acetyl-7-{4-[2-(2,4,5-trichlorophenoxy)ethoxy]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid (6-chlorobenzo[1,3]dioxol-5-ylmethyl)cyclopropylamide formate salt

30

Synthesized according to typical procedures H and E from bicyclononene BN9 and (6-chlorobenzo[1,3]dioxol-5-ylmethyl)cyclopropylamine. LC-MS: R_t = 0.94; ES+: 734.06.

Example 740

(*rac.*)-(1*R*, 5*S*)-7-{4-[3-(2,3,6-Trifluorophenoxy)propylphenyl]-3,9-diaza-bicyclo[3.3.1]non-6-ene-6-carboxylic acid (2-chlorobenzyl)cyclopropylamide

5

From Example 149, separated by chiral preparative HPLC.

Example 741

10 (*rac.*)-(1*S*, 5*R*)-7-{4-[3-(2,3,6-Trifluorophenoxy)propylphenyl]-3,9-diaza-bicyclo[3.3.1]non-6-ene-6-carboxylic acid (2-chlorobenzyl)cyclopropylamide

From Example 149, separated by chiral preparative HPLC.

15 **Example 742**

(*rac.*)-(1*R*^{*}, 5*S*^{*})-3-Acetyl-7-{4-[3-(2,6-dichloro-4-methylphenoxy)propyl]-phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid (2-chlorobenzyl)-cyclopropylamide formate salt

20

Synthesized according to typical procedures F and E from bicyclononene BB1 and 2,6-dichloro-4-methylphenol. LC-MS: R_t = 0.96; ES+: 666.35.

Example 743

25

(*rac.*)-(1*R*^{*}, 5*S*^{*})-7-{4-[3-(2-Bromo-5-fluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid (2-chlorobenzyl)cyclopropylamide

30

Synthesized according to typical procedure E from bicyclononene AL20. LC-MS: R_t = 0.86; ES+: 640.21.

The following assay was carried out in order to determine the activity of the compounds of general formula I and their salts.

5 **Inhibition of human recombinant renin by the compounds of the
invention**

10 The enzymatic in vitro assay was performed in 384-well polypropylene plates (Nunc). The assay buffer consisted of 10 mM PBS (Gibco BRL) including 1 mM EDTA and 0.1% BSA. The incubates were composed of 50 µL per well of an enzyme mix and 2.5 µL of renin inhibitors in DMSO. The enzyme mix was premixed at 4°C and consists of the following components:

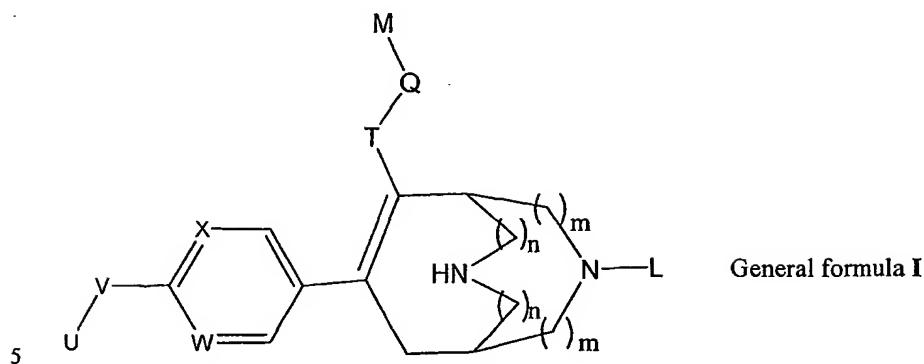
- human recombinant renin (0.16 ng/mL)
- synthetic human angiotensin(1-14) (0.5 µM)
- hydroxyquinoline sulfate (1 mM)

15 The mixtures were then incubated at 37°C for 3 h.

To determine the enzymatic activity and its inhibition, the accumulated Ang I was detected by an enzyme immunoassay (EIA) in 384-well plates (Nunc). 5 µL of the incubates or standards were transferred to immuno plates which were previously coated with a covalent complex of Ang I and bovine serum albumin (Ang I – BSA). 75 µL of Ang I-antibodies in essaybuffer above including 0.01% Tween 20 were added and a primary incubation made at 4 °C overnight. The plates were washed 3 times with PBS including 0.01% Tween 20, and then incubated for 2 h at rt with an antirabbit-peroxidase coupled antibody (WA 934, Amersham). After washing the plates 3 times, the *peroxidase substrate* ABTS (2,2'-azino-di-(3-ethylbenzthiazolinsulfonate), was added and the plates incubated for 60 min at room temperature. After stopping the reaction with 0.1 M citric acid pH 4.3 the plate was evaluated in a microplate reader at 405 nm. The percentage of inhibition was calculated of each concentration point and the concentration of renin inhibition was determined that inhibited the enzyme activity by 50% (IC_{50}). The IC_{50} -values of all compounds tested are below 100 nM. However selected compounds exhibit a very good bioavailability and are metabolically more stable than prior art compounds.

Claims

1. Compounds of the general formula I



wherein

X and W represent independently a nitrogen atom or a -CH- group;

10

V represents $-(\text{CH}_2)_r-$; $-\text{A}-(\text{CH}_2)_s-$; $-\text{CH}_2-\text{A}-(\text{CH}_2)_t-$; $-(\text{CH}_2)_s-\text{A}-$; $-(\text{CH}_2)_2-\text{A}-$
 $(\text{CH}_2)_u-$; $-\text{A}-(\text{CH}_2)_v-\text{B}-$; $-\text{CH}_2-\text{CH}_2-\text{CH}_2-\text{A}-\text{CH}_2-$; $-\text{A}-\text{CH}_2-\text{CH}_2-\text{B}-\text{CH}_2-$; $-\text{CH}_2-\text{A}-$
 $\text{CH}_2-\text{CH}_2-\text{B}-$; $-\text{CH}_2-\text{CH}_2-\text{CH}_2-\text{A}-\text{CH}_2-\text{CH}_2-$; $-\text{CH}_2-\text{CH}_2-\text{CH}_2-\text{CH}_2-\text{A}-\text{CH}_2-$; $-\text{A}-$
 $\text{CH}_2-\text{CH}_2-\text{B}-\text{CH}_2-\text{CH}_2-$; $-\text{CH}_2-\text{A}-\text{CH}_2-\text{CH}_2-\text{B}-\text{CH}_2-$; $-\text{CH}_2-\text{A}-\text{CH}_2-\text{CH}_2-\text{CH}_2-\text{B}-$; or
15 $-\text{CH}_2-\text{CH}_2-\text{A}-\text{CH}_2-\text{CH}_2-\text{B}-$;

A and B independently represent $-\text{O}-$; $-\text{S}-$; $-\text{SO}-$; $-\text{SO}_2-$;

U represents aryl; heteroaryl;

20

T represents $-\text{CONR}^1-$; $-(\text{CH}_2)_p\text{OCO}-$; $-(\text{CH}_2)_p\text{N}(\text{R}^1)\text{CO}-$; $-(\text{CH}_2)_p\text{N}(\text{R}^1)\text{SO}_2-$; or
 $-\text{COO}-$;

Q represents lower alkylene; lower alkenylene;

25

M represents hydrogen; cycloalkyl; aryl; heterocyclyl; heteroaryl;

L represents -R³; -COR³; -COOR³; -CONR²R³; -SO₂R³; -SO₂NR²R³;
-COCH(Aryl)₂;

5

R¹ represents hydrogen; lower alkyl; lower alkenyl; lower alkinyl; cycloalkyl;
aryl; cycloalkyl - lower alkyl;

10 R² and R^{2'} independently represent hydrogen; lower alkyl; lower alkenyl;
cycloalkyl; cycloalkyl - lower alkyl;

15 R³ represents hydrogen; lower alkyl; lower alkenyl; cycloalkyl; aryl; heteroaryl;
heterocyclyl; cycloalkyl - lower alkyl; aryl - lower alkyl; heteroaryl - lower alkyl;
heterocyclyl - lower alkyl; aryloxy - lower alkyl; heteroaryloxy - lower alkyl,
whereby these groups may be unsubstituted or mono-, di- or trisubstituted with
hydroxy, -OCOR², -COOR², lower alkoxy, cyano, -CONR²R^{2'}, -NH(NH)NH²,
-NR⁴R^{4'} or lower alkyl, with the proviso that a carbon atom is attached at the most
to one heteroatom in case this carbon atom is sp³-hybridized;

20 R⁴ and R^{4'} independently represents hydrogen; lower alkyl; cycloalkyl; cycloalkyl
- lower alkyl; hydroxy - lower alkyl; -COOR²; -CONH₂;

m and n represent the integer 0 or 1, with the proviso that in case m represents the
integer 1, n is the integer 0, and in case n represents the integer 1, m is the integer
25 0;

p is the integer 1, 2, 3 or 4;

r is the integer 3, 4, 5, or 6;

s is the integer 2, 3, 4, or 5;

30 t is the integer 1, 2, 3, or 4;

u is the integer 1, 2, or 3;

v is the integer 2, 3, or 4;

and optically pure enantiomers, mixtures of enantiomers such as racemates, diastereomers, mixtures of diastereomers, diastereomeric racemates, mixtures of diastereomeric racemates, and the meso-form; as well as pharmaceutically acceptable salts, solvent complexes and morphological forms.

2. Compounds of general formula I wherein X, W, V, U, T, Q, L, and M are as defined in general formula I and

10 n is 0

m is 1,

and optically pure enantiomers, mixtures of enantiomers such as racemates, diastereomers, mixtures of diastereomers, diastereomeric racemates, mixtures of diastereomeric racemates, and the meso-form; as well as pharmaceutically acceptable salts, solvent complexes and morphological forms.

3. Compounds of general formula I wherein X, W, V, U, T, Q, M, m, and n are as defined in general formula I and

20

L represents -COR^{3"}; -COOR^{3"}; -CONR^{2"}R^{3"};

R^{2"} and R^{3"} represent independently lower alkyl; lower cycloalkyl - lower alkyl, which lower alkyl and lower cycloalkyl-lower alkyl are undubstituted or mono-substituted with halogen, -CN, -OH, -OCOCH₃, -CONH₂, -COOH, or -NH₂, with the proviso that a carbon atom is attached at the most to one heteroatom in case this carbon atom is sp³-hybridized,

and optically pure enantiomers, mixtures of enantiomers such as racemates, diastereomers, mixtures of diastereomers, diastereomeric racemates, mixtures of diastereomeric racemates, and the meso-form; as well as pharmaceutically acceptable salts, solvent complexes and morphological forms.

4. Compounds of general formula I wherein X, W, V, U, L, m, and n are as defined in general formula I and

- 5 T represents -CONR¹;
Q represents methylene;
M represents hydrogen; aryl or heteroaryl;

and optically pure enantiomers, mixtures of enantiomers such as racemates,
10 diastereomers, mixtures of diastereomers, diastereomeric racemates, mixtures of diastereomeric racemates, and the meso-form; as well as pharmaceutically acceptable salts, solvent complexes and morphological forms.

5. Compounds of general formula I wherein X, W, U, L, T, Q, M, m, and n are as
15 defined in general formula I and

- V represents -CH₂CH₂O-; -CH₂CH₂CH₂O-; -OCH₂CH₂O-;

and optically pure enantiomers, mixtures of enantiomers such as racemates,
20 diastereomers, mixtures of diastereomers, diastereomeric racemates, mixtures of diastereomeric racemates, and the meso-form; as well as pharmaceutically acceptable salts, solvent complexes and morphological forms.

6. Compounds of general formula I wherein V, U, T, Q, M, L, m, and n are as
25 defined in general formula I and

- X and W represent a -CH- group

and optically pure enantiomers, mixtures of enantiomers such as racemates,
30 diastereomers, mixtures of diastereomers, diastereomeric racemates, mixtures of diastereomeric racemates, and the meso-form; as well as pharmaceutically acceptable salts, solvent complexes and morphological forms.

7. Compounds of general formula I wherein X, W, V, Q, T, M, L, m, and n are as defined in general formula I and

- 5 U is a mono-, di-, or trisubstituted phenyl whereby the substituents are halogen; lower alkyl or lower alkoxy

and optically pure enantiomers, mixtures of enantiomers such as racemates, diastereomers, mixtures of diastereomers, diastereomeric racemates, mixtures of 10 diastereomeric racemates, and the meso-form; as well as pharmaceutically acceptable salts, solvent complexes and morphological forms.

8. The compounds according to any one of claims 1 - 7 selected from the group consisting of

15

(*rac.*)-(2-methoxyphenyl)acetic acid (*IR**, *5S**)-7-{4-[3-(2-methoxybenzyloxy)-propoxy]phenyl}-3-(2-thiophen-2-ylacetyl)-3,9-diazabicyclo[3.3.1]non-6-en-6-ylmethyl ester;

20

(*rac.*)-(2-methoxyphenyl)acetic acid (*IR**, *5S**)-3-[2-(4-chlorophenyl)acetyl]-7-{4-[3-(2-methoxybenzyloxy)propoxy]phenyl}-3,9-diazabicyclo[3.3.1]non-6-en-6-ylmethyl ester;

25

(*rac.*)-(2-methoxyphenyl)acetic acid (*IR**, *5S**)-7-{4-[3-(2-methoxybenzyloxy)-propoxy]phenyl}-3-(quinoxaline-2-carbonyl)-3,9-diazabicyclo[3.3.1]non-6-en-6-ylmethyl ester;

30

(*rac.*)-(*IR**, *5S**)-3-acetyl-7-{4-[3-(2-methoxybenzyloxy)propoxy]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid [2-(2-chlorophenyl)ethyl]methylamide;

- (*rac.*)-(2-methoxyphenyl)acetic acid (*1R*^{*,} *5S*^{*})-3-(benzo[*b*]thiophene-3-carbonyl)-7-{4-[3-(2-methoxybenzyloxy)propoxy]phenyl}-3,9-diazabicyclo[3.3.1]non-6-en-6-ylmethyl ester;
- 5 (*rac.*)-(2-methoxyphenyl)acetic acid (*1R*^{*,} *5S*^{*})-3-acetyl-7-{4-[3-(2-methoxybenzyloxy)propoxy]phenyl}-3,9-diazabicyclo[3.3.1]non-6-en-6-ylmethyl ester;
- 10 (*rac.*)-(2-methoxyphenyl)acetic acid (*1R*^{*,} *5S*^{*})-7-{4-[3-(2-methoxybenzyloxy)propoxy]phenyl}-3-phenylmethanesulfonyl-3,9-diazabicyclo[3.3.1]non-6-en-6-ylmethyl ester;
- 15 (*rac.*)-(*1R*^{*,} *5S*^{*})-3-acetyl-7-{4-[3-(2-methoxybenzyloxy)propoxy]phenyl}-3,9-diaza-bicyclo[3.3.1]non-6-ene-6-carboxylic acid [2-(4-methoxyphenyl)ethyl]-methylamide;
- 20 (*rac.*)-(*1R*^{*,} *5S*^{*})-3-acetyl-7-{4-[2-(2-chloro-5-methylphenoxy)ethyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid methylphenethylamide;
- 25 (*rac.*)-(*1R*^{*,} *5S*^{*})-3-acetyl-7-{4-[2-(2-bromo-5-fluorophenoxy)ethyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid methylphenethylamide;
- 30 (*rac.*)-(2-methoxyphenyl)acetic acid (*1R*^{*,} *5S*^{*})-3-[2-(4-chlorophenyl)acetyl]-7-{6-[3-(2-methoxybenzyloxy)propoxy]pyridin-3-yl}-3,9-diazabicyclo[3.3.1]non-6-en-6-ylmethyl ester;
- 35 (*rac.*)-(*1R*^{*,} *5S*^{*})-3-acetyl-7-{4-[2-(2,5-dimethylphenoxy)ethyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid methylphenethylamide;
- 40 (*rac.*)-(*1R*^{*,} *5S*^{*})-3-acetyl-7-{4-[3-(2-methoxybenzyloxy)propoxy]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid [2-(4-chlorophenyl)ethyl]methylamide;

(*rac.*)-(1*R*^{*, 5*S*})-3-acetyl-7-{4-[3-(2-methoxybenzyloxy)propoxy]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid [2-(3-chlorophenyl)ethyl]methylamide;

5 (*rac.*)-(1*R*^{*, 5*S*})-3-acetyl-7-{4-[3-(2-methoxybenzyloxy)propoxy]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid ethylphenethylamide;

10 (*rac.*)-(1*R*^{*, 5*S*})-3-acetyl-7-{4-[3-(2-methoxybenzyloxy)propoxy]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid [2-(3-methoxyphenyl)ethyl]methylamide;

(*rac.*)-(1*R*^{*, 5*S*})-3-acetyl-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid methylphenethylamide;

15 (*rac.*)-(1*R*^{*, 5*S*})-3-acetyl-7-{4-[3-(2-methoxybenzyloxy)propoxy]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid methylphenethylamide;

(*rac.*)-(1*R*^{*, 5*S*})-3-acetyl-7-{4-[3-(2-bromo-5-fluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid methylphenethylamide;

20 (*rac.*)-(2-methoxyphenyl)acetic acid (1*R*^{*, 5*S*})-7-{4-[3-(2-methoxybenzyloxy)propoxy]phenyl}-3-methyl-3,9-diazabicyclo[3.3.1]non-6-en-6-ylmethyl ester;

25 (*rac.*)-(1*R*^{*, 5*S*})-3-acetyl-7-{4-[3-(2-methoxybenzyloxy)propoxy]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid [2-(3,4-dimethoxyphenyl)ethyl]methylamide;

(*rac.*)-(1*R*^{*, 5*S*})-3-acetyl-7-{4-[2-(2,3,5-trimethylphenoxy)ethyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid methylphenethylamide;

30 (*rac.*)-(1*R*^{*, 5*S*})-3-acetyl-7-{4-[3-(2,5-dichlorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid methylphenethylamide;

- (*rac.*)-(1*R*^{*}, 5*S*^{*})-3-acetyl-7-{4-[2-(2,3-dimethylphenoxy)ethyl]phenyl}-3,9-di-azabicyclo[3.3.1]non-6-ene-6-carboxylic acid methylphenethylamide;
- 5 (*rac.*)-*N*-(*IR*^{*}, 5*S*^{*})-3-acetyl-7-{4-[2-(2-chloro-4,5-dimethylphenoxy)ethyl]phe-nyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid methylphenethylamide;
- (*rac.*)-(2-methoxyphenyl)acetic acid (*IR*^{*}, 5*S*^{*})-7-{4-[3-(2-methoxybenzyloxy)-propoxy]phenyl}-3,9-diazabicyclo[3.3.1]non-6-en-6-ylmethyl ester;
- 10 (*rac.*)-*N*-(*IR*^{*}, 5*S*^{*})-3-acetyl-7-{4-[3-(2-methoxybenzyloxy)propoxy]phenyl}-3,9-diazabicyclo[3.3.1]non-6-en-6-ylmethyl)-2-(2-methoxyphenyl)-*N*-methyl-acetamide;
- 15 (*rac.*)-*N*-(*IR*^{*}, 5*S*^{*})-3-acetyl-7-{4-[2-(2-*tert*-butyl-4-methylphenoxy)ethyl]phe-nyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid methylphenethylamide;
- (*rac.*)-(1*R*^{*}, 5*S*^{*})-3-acetyl-7-{4-[2-(2,5-difluorophenoxy)ethyl]phenyl}-3,9-diaza-bicyclo[3.3.1]non-6-ene-6-carboxylic acid methylphenethylamide;
- 20 (*rac.*)-(1*R*^{*}, 5*S*^{*})-3-acetyl-7-{4-[3-(2-methoxybenzyloxy)propoxy]phenyl}-3,9-di-azabicyclo[3.3.1]non-6-ene-6-carboxylic acid methyl(3-phenylpropyl)amide;
- (*rac.*)-(1*R*^{*}, 5*S*^{*})-3-acetyl-7-{4-[3-(2,3-dichlorophenoxy)propyl]phenyl}-3,9-di-25 azabicyclo[3.3.1]non-6-ene-6-carboxylic acid methylphenethylamide;
- (*rac.*)-(1*R*^{*}, 5*S*^{*})-3-acetyl-7-{4-[3-(2-acetylphenoxy)propyl]phenyl}-3,9-diaza-bicyclo[3.3.1]non-6-ene-6-carboxylic acid methylphenethylamide;
- 30 (*rac.*)-(1*R*^{*}, 5*S*^{*})-3-acetyl-7-{4-[3-(2-methoxybenzyloxy)propoxy]phenyl}-3,9-di-azabicyclo[3.3.1]non-6-ene-6-carboxylic acid [2-(2-methoxyphenyl)ethyl]methyl-amide;

(*rac.*)-(1*R*^{*}, 5*S*^{*})-3-acetyl-7-{4-[3-(2-methoxybenzyloxy)propoxy]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid benzylmethyleamide;

5 (*rac.*)-(1*R*^{*}, 5*S*^{*})-3-acetyl-7-{4-[3-(2,5-difluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid methylphenethylamide;

(*rac.*)-(1*R*^{*}, 5*S*^{*})-3-acetyl-7-[4-(2-*o*-tolyloxyethyl)phenyl]-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid methylphenethylamide;

10 (*rac.*)-(1*R*^{*}, 5*S*^{*})-3-acetyl-7-{4-[2-(3-isopropylphenoxy)ethyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid methylphenethylamide;

15 (*rac.*)-(1*R*^{*}, 5*S*^{*})-3-acetyl-7-{4-[3-(2-chlorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid (2-chlorobenzyl)cyclopropylamide;

(*rac.*)-(1*R*^{*}, 5*S*^{*})-5-[7-{4-[2-(2-bromo-5-fluorophenoxy)ethyl]phenyl}-6-(methylphenethylcarbamoyl)-3,9-diazabicyclo[3.3.1]non-6-en-3-yl]-5-oxo-pentanoic acid;

20 (*rac.*)-(1*R*^{*}, 5*S*^{*})-3-acetyl-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid (2-chlorobenzyl)cyclopropylamide;

25 (*rac.*)-(1*R*^{*}, 5*S*^{*})-7-{4-[3-(2,3,6-trifluorophenoxy)propyl-phenyl]-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid (2-chlorobenzyl)cyclopropylamide;

(*rac.*)-5-((1*R*^{*}, 5*S*^{*})-6-[(2-chlorobenzyl)cyclopropylcarbamoyl]-7-{4-[3-(2,3,6-trifluorophenoxy)propylphenyl]-3,9-diazabicyclo[3.3.1]non-6-en-3-yl}-5-oxo-pentanoic acid;

- (*rac.*)-(1*R*^{*}, 5*S*^{*})-6-[(2-chlorobenzyl)cyclopropylcarbamoyl]-7-{4-[3-(2,3,6-trifluoro-phenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-9-carboxylic acid 2,2,2-trichloro-1,1-dimethylethyl ester;
- 5 (*rac.*)-5-((1*R*^{*}, 5*S*^{*})-6-[(2-chlorobenzyl)cyclopropylcarbamoyl]-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-en-3-yl)-2,2-dimethyl-5-oxo-pentanoic acid;
- 10 (*rac.*)-5-((1*R*^{*}, 5*S*^{*})-6-[(2-chlorobenzyl)cyclopropylcarbamoyl]-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-en-3-yl)-5-oxo-pentanoic acid methyl ester;
- 15 1:1-Mixture of (1*R*, 5*S*)-3-((1*S*, 4*R*)-4-hydroxypyrrolidine-2-carbonyl)-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid (2-chlorobenzyl)cyclopropylamide and (1*S*, 5*R*)-3-((1*S*, 4*R*)-4-hydroxypyrrolidine-2-carbonyl)-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid (2-chlorobenzyl)cyclopropylamide;
- 20 20 (*rac.*)-(1*R*^{*}, 5*S*^{*})-3-(4-carbamoylbutyryl)-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid (2-chlorobenzyl)cyclopropylamide;
- 25 (*rac.*)-(1*R*^{*}, 5*S*^{*})-3-(4-carbamoylbutyryl)-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid benzylcyclopropylamide;
- 30 (*rac.*)-(1*R*^{*}, 5*S*^{*})-3-(4-carbamoylbutyryl)-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid (2-chlorobenzyl)ethylamide;

- (*rac.*)-(1*R*^{*}, 5*S*^{*})-3-(4-carbamoylbutyryl)-7-{4-[3-(2,3,6-trifluorophenoxy)-propyl]phenyl}-3,9-diaza-bicyclo[3.3.1]non-6-ene-6-carboxylic acid cyclopropyl-(2-fluorobenzyl)amide;
- 5 (*rac.*)-(1*R*^{*}, 5*S*^{*})-3-(4-carbamoylbutyryl)-7-{4-[3-(2,3,6-trifluorophenoxy)-propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid cyclopropyl-(3-trifluoromethylbenzyl)amide;
- 10 (*rac.*)-(1*R*^{*}, 5*S*^{*})-3-(4-carbamoylbutyryl)-7-{4-[3-(2,3,6-trifluorophenoxy)-propyl]phenyl}-3,9-diaza-bicyclo[3.3.1]non-6-ene-6-carboxylic acid cyclopropyl-(2-methylbenzyl)amide;
- 15 (*rac.*)-(1*R*^{*}, 5*S*^{*})-3-(4-carbamoylbutyryl)-7-{4-[3-(2,3,6-trifluorophenoxy)-propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid cyclopropyl-phenethylamide;
- 20 (*rac.*)-(1*R*^{*}, 5*S*^{*})-3-(4-carbamoylbutyryl)-7-{4-[3-(2,3,6-trifluorophenoxy)-propyl]phenyl}-3,9-diaza-bicyclo[3.3.1]non-6-ene-6-carboxylic acid [2-(2-chlorophenyl)ethyl]cyclopropylamide;
- 25 (*rac.*)-(1*R*^{*}, 5*S*^{*})-3-(4-carbamoylbutyryl)-7-{4-[3-(2,3,6-trifluorophenoxy)-propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid cyclopropyl-(3,5-dimethoxybenzyl)amide;
- 30 1:1-mixture of (*rac.*)-(3*R*^{*})-5-((1*R*^{*}, 5*S*^{*})-6-[(2-chlorobenzyl)cyclopropylcarbamoyl]-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo-[3.3.1]non-6-en-3-yl)-3-hydroxy-5-oxopentanoic acid and (*rac.*)-(3*S*^{*})-5-((1*R*^{*}, 5*S*^{*})-6-[(2-chlorobenzyl)cyclopropyl-carbamoyl]-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo-[3.3.1]non-6-en-3-yl)-3-hydroxy-5-oxopentanoic acid;

- 1:1-mixture of (*rac.*)-(3*R**)-5-((1*R**, 5*S**)-6-(benzylcyclopropylcarbamoyl)-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-en-3-yl)-3-hydroxy-5-oxopentanoic acid and (*rac.*)-(3*S**)-5-((1*R**, 5*S**)-6-(benzylcyclopropylcarbamoyl)-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-en-3-yl)-3-hydroxy-5-oxopentanoic acid;
- 5
- 1:1-mixture of (*rac.*)-(3*R**)-5-((1*R**, 5*S**)-6-[(2-chlorobenzyl)ethylcarbamoyl]-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-en-3-yl)-3-hydroxy-5-oxopentanoic acid and (*rac.*)-(3*S**)-5-((1*R**, 5*S**)-6-[(2-chlorobenzyl)ethylcarbamoyl]-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-en-3-yl)-3-hydroxy-5-oxopentanoic acid;
- 10
- 1:1-mixture of (*rac.*)-(3*R**)-5-((1*R**, 5*S**)-6-[cyclopropyl-(2-fluorobenzyl)-carbamoyl]-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-en-3-yl)-3-hydroxy-5-oxopentanoic acid and (*rac.*)-(3*S**)-5-((1*R**, 5*S**)-6-[cyclopropyl-(2-fluorobenzyl)-carbamoyl]-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-en-3-yl)-3-hydroxy-5-oxopentanoic acid;
- 15
- 20 1:1-mixture of (*rac.*)-(3*R**)-5-((1*R**, 5*S**)-6-[cyclopropyl-(3-trifluoromethylbenzyl)carbamoyl]-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]-phenyl}-3,9-diazabicyclo[3.3.1]non-6-en-3-yl)-3-hydroxy-5-oxopentanoic acid and (*rac.*)-(3*S**)-5-((1*R**, 5*S**)-6-[cyclopropyl-(3-trifluoromethylbenzyl)carbamoyl]-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]-phenyl}-3,9-diazabicyclo[3.3.1]non-6-en-3-yl)-3-hydroxy-5-oxopentanoic acid;
- 25
- 30 1:1-mixture of (*rac.*)-(3*R**)-5-((1*R**, 5*S**)-6-[cyclopropyl-(2-methylbenzyl)-carbamoyl]-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-en-3-yl)-3-hydroxy-5-oxopentanoic acid and (*rac.*)-(3*S**)-5-((1*R**, 5*S**)-6-[cyclopropyl-(2-methylbenzyl)carbamoyl]-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-en-3-yl)-3-hydroxy-5-oxopentanoic acid;

- 1:1-mixture of (*rac.*)-(3*R**)-5-((1*R**, 5*S**)-6-{cyclopropyl-[2-(4-methoxyphenoxy)ethyl]carbamoyl}-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-en-3-yl)-3-hydroxy-5-oxopentanoic acid and (*rac.*)-(3*S**)-5-((1*R**, 5*S**)-6-{cyclopropyl-[2-(4-methoxyphenoxy)ethyl]carbamoyl}-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-en-3-yl)-3-hydroxy-5-oxopentanoic acid;
- 1:1-mixture of (*rac.*)-(3*R**)-5-((1*R**, 5*S**)-6-[cyclopropyl-(2-*m*-tolyloxyethyl)-carbamoyl]-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-en-3-yl)-3-hydroxy-5-oxopentanoic acid and (*rac.*)-(3*S**)-5-((1*R**, 5*S**)-6-[cyclopropyl-(2-*m*-tolyloxyethyl)-carbamoyl]-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-en-3-yl)-3-hydroxy-5-oxopentanoic acid;
- 15 1:1-mixture of (*rac.*)-(3*R**)-5-((1*R**, 5*S**)-6-{[2-(2-chlorophenyl)ethyl]-cyclopropylcarbamoyl}-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-en-3-yl)-3-hydroxy-5-oxopentanoic acid and (*rac.*)-(3*S**)-5-((1*R**, 5*S**)-6-{[2-(2-chlorophenyl)ethyl]cyclopropylcarbamoyl}-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-en-3-yl)-3-hydroxy-5-oxopentanoic acid;
- 20 1:1-mixture of (*rac.*)-(3*R**)-5-((1*R**, 5*S**)-6-{cyclopropyl-(3,5-dimethoxybenzyl)carbamoyl}-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-en-3-yl)-3-hydroxy-5-oxopentanoic acid and (*rac.*)-(3*S**)-5-((1*R**, 5*S**)-6-[cyclopropyl-(3,5-dimethoxybenzyl)carbamoyl]-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-en-3-yl)-3-hydroxy-5-oxopentanoic acid;
- 25 30 (*rac.*)-(1*R**, 5*S**)-3-acetyl-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid (2-chlorobenzyl)ethylamide;

- (*rac.*)-(1*R*^{*}, 5*S*^{*})-3-acetyl-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid cyclopropyl(3,5-dimethoxybenzyl)amide;
- 5 (*rac.*)-5-((1*R*^{*}, 5*S*^{*})-6-(benzylcyclopropylcarbamoyl)-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-en-3-yl)-5-oxopentanoic acid;
- 10 (*rac.*)-5-((1*R*^{*}, 5*S*^{*})-6-[(2-chlorobenzyl)ethylcarbamoyl]-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-en-3-yl)-5-oxopentanoic acid;
- 15 (*rac.*)-5-((1*R*^{*}, 5*S*^{*})-6-[cyclopropyl-(2-fluorobenzyl)carbamoyl]-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-en-3-yl)-5-oxopentanoic acid;
- 20 (*rac.*)-5-((1*R*^{*}, 5*S*^{*})-6-[cyclopropyl(3-trifluoromethylbenzyl)carbamoyl]-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-en-3-yl)-5-oxopentanoic acid;
- 25 (*rac.*)-5-((1*R*^{*}, 5*S*^{*})-6-{[2-(2-chlorophenyl)ethyl]cyclopropylcarbamoyl}-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-en-3-yl)-5-oxopentanoic acid;
- 30 (*rac.*)-5-((1*R*^{*}, 5*S*^{*})-6-[cyclopropyl(3,5-dimethoxybenzyl)carbamoyl]-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-en-3-yl)-5-oxopentanoic acid;

- (*rac.*.)-5-((*1R*^{*}, *5S*^{*})-6-[cyclopropyl(2-*p*-tolylethyl)carbamoyl]-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-en-3-yl)-5-oxopentanoic acid;
- 5 (*rac.*.)-5-((*1R*^{*}, *5S*^{*})-6-[(2-chlorobenzyl)ethylcarbamoyl]-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-en-3-yl)-5-oxopentanoic acid methyl ester;
- 10 (*rac.*.)-5-((*1R*^{*}, *5S*^{*})-6-[cyclopropyl-(2-fluorobenzyl)carbamoyl]-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-en-3-yl)-5-oxopentanoic acid methyl ester;
- 15 (*rac.*.)-5-((*1R*^{*}, *5S*^{*})-6-[cyclopropyl-(3,5-dimethoxybenzyl)carbamoyl]-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-en-3-yl)-5-oxopentanoic acid methyl ester;
- 20 (*rac.*.)-5-((*1R*^{*}, *5S*^{*})-6-[(2-chlorobenzyl)ethylcarbamoyl]-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-en-3-yl)-2,2-dimethyl-5-oxopentanoic acid;
- 25 (*rac.*.)-5-((*1R*^{*}, *5S*^{*})-6-[cyclopropyl-(2-fluorobenzyl)carbamoyl]-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-en-3-yl)-2,2-dimethyl-5-oxopentanoic acid;
- 30 (*rac.*.)-5-((*1R*^{*}, *5S*^{*})-6-[cyclopropyl-(3,5-dimethoxybenzyl)carbamoyl]-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-en-3-yl)-2,2-dimethyl-5-oxopentanoic acid;

- 1:1-mixture of (*rac.*)-(2*R*^{*, 3*S*^{*})-4-((1*R*^{*, 5*S*^{*})-6-[cyclopropyl-(2-methylbenzyl)-carbamoyl]-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-en-3-yl)-2,3-dihydroxy-4-oxobutyric acid and (*rac.*)-(2*S*^{*, 3*R*^{*})-4-((1*R*^{*, 5*S*^{*})-6-[cyclopropyl-(2-methylbenzyl)carbamoyl]-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-en-3-yl)-2,3-dihydroxy-4-oxobutyric acid;}}}}
- 5
10 (*rac.*)-5-{(1*R*^{*, 5*S*^{*})-7-{4-[3-(2-bromo-5-fluorophenoxy)propyl]phenyl}-6-[cyclopropyl-(2-fluorobenzyl)carbamoyl]-3,9-diazabicyclo[3.3.1]non-6-en-3-yl}-5-oxopentanoic acid;}
- 15
15 (*rac.*)-5-{(1*R*^{*, 5*S*^{*})-7-{4-[3-(2-bromo-5-fluorophenoxy)propyl]phenyl}-6-[cyclopropyl-[2-(3-methylphenoxy)ethyl]carbamoyl]-3,9-diazabicyclo[3.3.1]-non-6-en-3-yl)-5-oxopentanoic acid;}
- 20
20 (*rac.*)-5-((1*R*^{*, 5*S*^{*})-7-{4-[3-(2-bromo-5-fluorophenoxy)propyl]phenyl}-6-{[2-(2-chlorophenyl)ethyl]cyclopropylcarbamoyl}-3,9-diazabicyclo[3.3.1]non-6-en-3-yl)-5-oxopentanoic acid;}
- 25
25 (*rac.*)-5-((1*R*^{*, 5*S*^{*})-7-{4-[3-(2-bromo-5-fluorophenoxy)propyl]phenyl}-6-{[2-(2,3-difluorophenyl)ethyl]cyclopropylcarbamoyl}-3,9-diazabicyclo[3.3.1]non-6-en-3-yl)-5-oxopentanoic acid;}
- 30
30 (*rac.*)-5-((1*R*^{*, 5*S*^{*})-7-{4-[3-(2-bromo-5-fluorophenoxy)propyl]phenyl}-6-{[2-(2-methylphenyl)ethyl]cyclopropylcarbamoyl}-3,9-diazabicyclo[3.3.1]non-6-en-3-yl)-5-oxopentanoic acid;}

(*rac.*.)-5-((*1R*^{*}, *5S*^{*})-7-{4-[3-(2-bromo-5-fluorophenoxy)propyl]phenyl}-6-[[2-(4-methylphenyl)ethyl]cyclopropylcarbamoyl]-3,9-diazabicyclo[3.3.1]non-6-en-3-yl)-5-oxopentanoic acid;

5 (*rac.*.)-5-((*1R*^{*}, *5S*^{*})-7-{4-[3-(2-bromo-5-fluorophenoxy)propyl]phenyl}-6-[cyclopropyl-(3,5-dimethoxybenzyl)carbamoyl]-3,9-diazabicyclo[3.3.1]non-6-en-3-yl)-5-oxopentanoic acid;

10 (*rac.*.)-5-((*1R*^{*}, *5S*^{*})-7-{4-[3-(2-bromo-5-fluorophenoxy)propyl]phenyl}-6-[cyclopropyl-(3,5-dimethoxybenzyl)carbamoyl]-3,9-diazabicyclo[3.3.1]non-6-en-3-yl)-5-oxopentanoic acid methyl ester;

15 (*rac.*.)-5-((*1R*^{*}, *5S*^{*})-7-{4-[3-(2-bromo-5-fluorophenoxy)propyl]phenyl}-6-[cyclopropyl-(2-fluorobenzyl)carbamoyl]-3,9-diazabicyclo[3.3.1]non-6-en-3-yl)-2,2-dimethyl-5-oxo-pentanoic acid;

20 (*rac.*.)-5-((*1R*^{*}, *5S*^{*})-7-{4-[3-(2-bromo-5-fluorophenoxy)propyl]phenyl}-6-[cyclopropyl-(2-methylbenzyl)carbamoyl]-3,9-diazabicyclo[3.3.1]non-6-en-3-yl)-2,2-dimethyl-5-oxo-pentanoic acid;

25 (*rac.*.)-5-[(*1R*^{*}, *5S*^{*})-7-{4-[3-(2-bromo-5-fluorophenoxy)propyl]phenyl}-6-(cyclopropylphenethylcarbamoyl)-3,9-diazabicyclo[3.3.1]non-6-en-3-yl]-2,2-dimethyl-5-oxopentanoic acid;

30 (*rac.*.)-5-((*1R*^{*}, *5S*^{*})-7-{4-[3-(2-bromo-5-fluorophenoxy)propyl]phenyl}-6-[[2-(4-methylphenyl)ethyl]cyclopropylcarbamoyl]-3,9-diazabicyclo[3.3.1]non-6-en-3-yl)-2,2-dimethyl-5-oxopentanoic acid;

35 (*rac.*.)-5-((*1R*^{*}, *5S*^{*})-7-{4-[3-(2-bromo-5-fluorophenoxy)propyl]phenyl}-6-[cyclopropyl-(3,5-dimethoxybenzyl)carbamoyl]-3,9-diazabicyclo[3.3.1]non-6-en-3-yl)-2,2-dimethyl-5-oxopentanoic acid;

- 1:1-mixture of (*rac.*)-(2*R*^{*}, 3*S*^{*})-4-{(1*R*^{*}, 5*S*^{*})-7-{4-[3-(2-bromo-5-fluoro-phenoxy)propyl]phenyl}-6-[cyclopropyl-(3,5-dimethoxybenzyl)carbamoyl]-3,9-diazabicyclo[3.3.1]non-6-en-3-yl}-2,3-dihydroxy-4-oxobutyric acid and (*rac.*)-(2*S*^{*}, 3*R*^{*})-4-{(1*R*^{*}, 5*S*^{*})-7-{4-[3-(2-bromo-5-fluoro-phenoxy)propyl]phenyl}-6-[cyclopropyl-(3,5-dimethoxybenzyl)carbamoyl]-3,9-diazabicyclo[3.3.1]non-6-en-3-yl}-2,3-dihydroxy-4-oxobutyric acid;
- 5 (10) (*rac.*)-5-((1*R*^{*}, 5*S*^{*})-6-[(2-chlorobenzyl)cyclopropylcarbamoyl]-7-{4-[2-(2,3,5-trimethylphenoxy)ethyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-en-3-yl)-5-oxo-pentanoic acid;
- 15 (15) (*rac.*)-5-((1*R*^{*}, 5*S*^{*})-6-[(3-trifluoromethylbenzyl)cyclopropylcarbamoyl]-7-{4-[2-(2,3,5-trimethylphenoxy)ethyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-en-3-yl)-5-oxo-pentanoic acid;
- 20 (20) (*rac.*)-5-((1*R*^{*}, 5*S*^{*})-6-[(2-methylbenzyl)cyclopropylcarbamoyl]-7-{4-[2-(2,3,5-trimethylphenoxy)ethyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-en-3-yl)-5-oxo-pentanoic acid;
- 25 (25) (*rac.*)-5-((1*R*^{*}, 5*S*^{*})-6-[(2-chlorobenzyl)ethylcarbamoyl]-7-{4-[2-(2,3,5-trimethylphenoxy)ethyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-en-3-yl)-2,2-dimethyl-5-oxopentanoic acid;
- 30 (30) (*rac.*)-5-((1*R*^{*}, 5*S*^{*})-7-{4-[3-(2-bromo-5-fluorophenoxy)propyl]phenyl}-6-[(2-chlorobenzyl)cyclopropylcarbamoyl]-3,9-diazabicyclo[3.3.1]non-6-en-3-yl)-5-oxopentanoic acid;

- (*rac.*.)-5-{(1*R**, 5*S**)-7-{4-[3-(2-bromo-5-fluorophenoxy)propyl]phenyl}-6-[(2-chlorobenzyl)cyclopropylcarbamoyl]-3,9-diazabicyclo[3.3.1]non-6-en-3-yl}-2,2-dimethyl-5-oxopentanoic acid;
- 5 (*rac.*.)-5-{(1*R**, 5*S**)-4-[3-(2-bromo-5-fluorophenoxy)propyl]phenyl}-3-(4-carbamoylbutyryl)-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid (2-chlorobenzyl)cyclopropylamide;
- 10 (*rac.*.)-5-((1*R**, 5*S**)-6-(benzylcyclopropylcarbamoyl)-7-{4-[3-(2-bromo-5-fluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-en-3-yl)-5-oxopentanoic acid methyl ester;
- 15 (*rac.*.)-(1*R**, 5*S**)-3-acetyl-7-{4-[2-(2-chloro-4,5-dimethylphenoxy)ethoxy]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid (2-chlorobenzyl)-cyclopropylamide;
- 20 (*rac.*.)-5-((1*R**, 5*S**)-6-[(2-chlorobenzyl)cyclopropylcarbamoyl]-7-{4-[2-(2-chloro-4,5-dimethylphenoxy)ethoxy]phenyl}-3,9-diazabicyclo[3.3.1]non-6-en-3-yl)-5-oxopentanoic acid;
- 25 (*rac.*.)-5-{(1*R**, 5*S**)-7-{4-[2-(2-chloro-4,5-dimethylphenoxy)ethoxy]phenyl}-6-[cyclopropyl-(2-methylbenzyl)carbamoyl]-3,9-diazabicyclo[3.3.1]non-6-en-3-yl}-5-oxopentanoic acid;
- 30 (*rac.*.)-5-{(1*R**, 5*S**)-7-{4-[2-(2-chloro-4,5-dimethylphenoxy)ethoxy]phenyl}-6-[cyclopropyl-(3,5-dimethoxybenzyl)carbamoyl]-3,9-diazabicyclo[3.3.1]non-6-en-3-yl}-5-oxopentanoic acid;
- (*rac.*.)-5-((1*R**, 5*S**)-6-[(2-chlorobenzyl)cyclopropylcarbamoyl]-7-{4-[2-(2-chloro-4,5-dimethylphenoxy)ethoxy]phenyl}-3,9-diazabicyclo[3.3.1]non-6-en-3-yl)-2,2-dimethyl-5-oxopentanoic acid;

(*rac.*.)-5-((*1R*^{*}, *5S*^{*})-6-[(2-chlorobenzyl)ethylcarbamoyl]-7-{4-[2-(2-chloro-4,5-dimethylphenoxy)ethoxy]phenyl}-3,9-diazabicyclo[3.3.1]non-6-en-3-yl)-2,2-dimethyl-5-oxopentanoic acid;

5 (rac.)-5-((*1R*^{*}, *5S*^{*})-7-{4-[2-(2-chloro-4,5-dimethylphenoxy)ethoxy]phenyl}-6-[cyclopropyl-(2-fluorobenzyl)carbamoyl]-3,9-diazabicyclo[3.3.1]non-6-en-3-yl)-2,2-dimethyl-5-oxopentanoic acid;

10 (rac.)-5-((*1R*^{*}, *5S*^{*})-6-[cyclopropyl-(3,5-dimethoxybenzyl)carbamoyl]-7-{4-[2-(2,3,5-trimethylphenoxy)ethyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-en-3-yl)-5-oxopentanoic acid;

15 (rac.)-5-((*1R*^{*}, *5S*^{*})-6-[cyclopropyl-(2-*p*-tolylethyl)carbamoyl]-7-{4-[2-(2,3,5-trimethylphenoxy)ethyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-en-3-yl)-5-oxopentanoic acid;

20 (rac.)-5-((*1R*^{*}, *5S*^{*})-7-{4-[2-(2-chloro-4,5-dimethylphenoxy)ethoxy]phenyl}-6-[cyclopropyl-[2-(2-hydroxyethyl)benzyl]carbamoyl]-3,9-diazabicyclo[3.3.1]non-6-en-3-yl)-5-oxopentanoic acid;

25 (rac.)-5-((*1R*^{*}, *5S*^{*})-6-(cyclopropylphenethylcarbamoyl)-7-{4-[2-(2,3,5-trimethylphenoxy)ethyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-en-3-yl)-2,2-dimethyl-5-oxopentanoic acid;

25 (rac.)-((*1R*^{*}, *5S*^{*})-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid (2-chlorobenzyl)ethylamide;

30 (rac.)-((*1R*^{*}, *5S*^{*})-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-3,6-dicarboxylic acid 6-[(2-chlorobenzyl)cyclopropyl-
amide] 3-dimethylamide;

- (*rac.*)-(1*R*^{*}, 5*S*^{*})-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-3,6-dicarboxylic acid 6-[(2-chlorobenzyl)cyclopropylamide] 3-diethylamide;
- 5 (*rac.*)-(1*R*^{*}, 5*S*^{*})-6-[(2-chlorobenzyl)cyclopropylcarbamoyl]-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-3-carboxylic acid methyl ester;
- 10 (*rac.*)-(1*R*^{*}, 5*S*^{*})-6-[(2-chlorobenzyl)cyclopropylcarbamoyl]-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-3-carboxylic acid ethyl ester;
- 15 (*rac.*)-(1*R*^{*}, 5*S*^{*})-3-methanesulfonyl-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid (2-chlorobenzyl)-cyclopropylamide;
- 20 (*rac.*)-(1*R*^{*}, 5*S*^{*})-3-ethanesulfonyl-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid (2-chlorobenzyl)-cyclopropylamide;
- 25 (*rac.*)-5-((1*R*^{*}, 5*S*^{*})-6-[(2-chlorobenzyl)cyclopropylcarbamoyl]-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-en-3-yl)-5-oxo-pentanoic acid ethyl ester;
- 30 (*rac.*)-4-((1*R*^{*}, 5*S*^{*})-6-[(2-chlorobenzyl)cyclopropylcarbamoyl]-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-en-3-yl)-4-oxo-butrylic acid;
- 35 (*rac.*)-3-[((1*R*^{*}, 5*S*^{*})-6-[(2-chlorobenzyl)cyclopropylcarbamoyl]-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-3-carbonyl)-amino]propionic acid ethyl ester;

(*rac.*)4-[((*1R*^{*}, *5S*^{*})-6-[(2-chlorobenzyl)cyclopropylcarbamoyl]-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-3-carbonyl)-amino]butyric acid ethyl ester;

5 (*rac.*)-(*1R*^{*}, *5S*^{*})-3-(3-carbamoylpropionyl)-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid (2-chlorobenzyl)cyclopropylamide;

10 (*rac.*)-(*1R*^{*}, *5S*^{*})-3-(2-hydroxyacetyl)-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid (2-chlorobenzyl)cyclopropylamide;

15 (*1S*, *5R*)-3-((*3R*)-3-hydroxybutyryl)-7-{4-[3-(2,3,6-trifluorophenoxy)-propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid (2-chlorobenzyl)cyclopropylamide;

20 1:1-mixture of (*rac.*)-(*1R*^{*}, *5S*^{*})-3-((*1R*^{*}, *2S*^{*})-2-hydroxycyclopentanecarbonyl)-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid (2-chlorobenzyl)cyclopropylamide and (*rac.*)-(*1R*^{*}, *5S*^{*})-3-((*1S*^{*}, *2R*^{*})-2-hydroxycyclopentanecarbonyl)-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid (2-chlorobenzyl)cyclopropylamide;

25 (*rac.*)-(*1R*^{*}, *5S*^{*})-9-acetyl-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid (2-chlorobenzyl)cyclopropylamide;

30 (*rac.*)-5-((*1R*^{*}, *5S*^{*})-6-[(2-chlorobenzyl)cyclopropylcarbamoyl]-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-en-9-yl)-5-oxo-pentanoic acid;

(*rac.*.)-5-((*1R**, *5S**)-6-[(2-chlorobenzyl)cyclopropylcarbamoyl]-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-en-9-yl)-5-oxo-pentanoic acid ethyl ester;

5 (*rac.*.)-5-((*1R**, *5S**)-6-[(2-chlorobenzyl)cyclopropylcarbamoyl]-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-en-9-yl)-5-oxo-pentanoic acid methyl ester;

10 1:1-mixture of (*rac.*.)-(*3R**)-5-((*1R**, *5S**)-6-[(2-chlorobenzyl)cyclopropylcarbamoyl]-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-en-9-yl)-3-hydroxy-5-oxo-pentanoic acid and (*rac.*.)-(*3S**)-5-((*1R**, *5S**)-6-[(2-chlorobenzyl)cyclopropylcarbamoyl]-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-en-9-yl)-3-hydroxy-5-oxo-pentanoic acid;

15 (*rac.*.)-5-((*1R**, *5S**)-6-[(2-chlorobenzyl)cyclopropylcarbamoyl]-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-en-9-yl)-2,2-dimethyl-5-oxopentanoic acid;

20 (*rac.*.)-4-((*1R**, *5S**)-6-[(2-chlorobenzyl)cyclopropylcarbamoyl]-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-en-9-yl)-4-oxo-butyric acid;

25 (*rac.*.)-((*1R**, *5S**)-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6,9-dicarboxylic acid 6-[(2-chlorobenzyl)cyclopropylamide] 9-dimethylamide;

30 (*rac.*.)-((*1R**, *5S**)-6-[(2-chlorobenzyl)cyclopropylcarbamoyl]-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-9-carboxylic acid methyl ester;

- (*rac.*)-(1*R*^{*}, 5*S*^{*})-6-[(2-chlorobenzyl)cyclopropylcarbamoyl]-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-9-carboxylic acid ethyl ester;
- 5 (*rac.*)-3-[(1*R*^{*}, 5*S*^{*})-(6-[(2-chlorobenzyl)cyclopropylcarbamoyl]-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-9-carboxyl)-amino]propionic acid ethyl ester;
- 10 (*rac.*)-4-[((1*R*^{*}, 5*S*^{*})-6-[(2-chlorobenzyl)cyclopropylcarbamoyl]-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-9-carboxyl)-amino]butyric acid ethyl ester;
- 15 (*rac.*)-(1*R*^{*}, 5*S*^{*})-3-formyl-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid (2-chlorobenzyl)cyclopropylamide;
- 20 (*rac.*)-3-[((1*R*^{*}, 5*S*^{*})-6-[(2-chlorobenzyl)cyclopropylcarbamoyl]-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-3-carboxyl)-amino]propionic acid;
- 25 (*rac.*)-4-[((1*R*^{*}, 5*S*^{*})-6-[(2-chlorobenzyl)cyclopropylcarbamoyl]-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-9-carboxyl)-amino]butyric acid;
- 30 (*rac.*)-(1*R*^{*}, 5*S*^{*})-3-acetyl-7-{4-[2-(2,3,6-trifluorophenoxy)ethyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid (2-chlorobenzyl)cyclopropylamide;

- (*rac.*)-(1*R**, 5*S**)-9-(4-carbamoylbutyryl)-7-{4-[2-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diaza-bicyclo[3.3.1]non-6-ene-6-carboxylic acid (2-chlorobenzyl)cyclopropylamide;
- 5 (*rac.*)-(1*R**, 5*S**)-9-(3-carbamoylpropionyl)-7-{4-[2-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diaza-bicyclo[3.3.1]non-6-ene-6-carboxylic acid (2-chlorobenzyl)cyclopropylamide;
- 10 (*rac.*)-(1*R**, 5*S**)-9-(2-hydroxyacetyl)-7-{4-[2-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diaza-bicyclo[3.3.1]non-6-ene-6-carboxylic acid (2-chlorobenzyl)cyclopropylamide;
- 15 (1*R*, 5*S*)-9-((3*S*)-3-hydroxybutyryl)-7-{4-[2-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid (2-chlorobenzyl)cyclopropylamide;
- 20 (*rac.*)-(1*R**, 5*S**)-9-methanesulfonyl-7-{4-[2-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid (2-chlorobenzyl)cyclopropylamide;
- 25 (*rac.*)-(1*R**, 5*S**)-9-ethanesulfonyl-7-{4-[2-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid (2-chlorobenzyl)cyclopropylamide;
- 30 (*rac.*)-(1*R**, 5*S**)-3-acetyl-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid cyclopropyl-(2-fluoro-5-methoxybenzyl)amide;
- 35 (*rac.*)-(1*R**, 5*S**)-3-acetyl-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid cyclopropyl-(3-methoxybenzyl)amide;

(*rac.*)-(1*R*^{*}, 5*S*^{*})-3-acetyl-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid cyclopropyl-(3,4-dimethoxybenzyl)amide;

5 (*rac.*)-(1*R*^{*}, 5*S*^{*})-3-acetyl-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid (2-chloro-3-trifluoromethylbenzyl)cyclopropylamide;

10 (*rac.*)-(1*R*^{*}, 5*S*^{*})-3-acetyl-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid benzo[1,3]dioxol-5-ylmethylcyclopropylamide;

15 (*rac.*)-(1*R*^{*}, 5*S*^{*})-3-acetyl-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid (2-chloro-6-fluorobenzyl)cyclopropylamide;

20 (*rac.*)-(1*R*^{*}, 5*S*^{*})-3-acetyl-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid (2-bromobenzyl)cyclopropylamide;

25 (*rac.*)-(1*R*^{*}, 5*S*^{*})-3-acetyl-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid cyclopropyl-(2,3-dimethylbenzyl)amide;

30 (*rac.*)-(1*R*^{*}, 5*S*^{*})-3-acetyl-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid cyclopropyl-(3,5-difluorobenzyl)amide;

35 (*rac.*)-(1*R*^{*}, 5*S*^{*})-3-acetyl-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid cyclopropyl-(2,3-dichlorobenzyl)amide;

- (*rac.*)-(1*R*^{*}, 5*S*^{*})-3-acetyl-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid cyclopropyl-(3-trifluoromethoxybenzyl)amide;
- 5 (*rac.*)-(1*R*^{*}, 5*S*^{*})-3-acetyl-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid cyclopropyl-(3-methylbenzyl)amide;
- 10 (*rac.*)-(1*R*^{*}, 5*S*^{*})-3-acetyl-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diaza-bicyclo[3.3.1]non-6-ene-6-carboxylic acid (3-chlorobenzyl)cyclopropylamide;
- 15 (*rac.*)-(1*R*^{*}, 5*S*^{*})-7-{4-[3-(2-bromo-5-fluorophenoxy)propyl]phenyl}-3-(4-carbamoylbutyryl)-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid cyclopropyl-(3,5-dimethoxybenzyl)amide;
- 20 (*rac.*)-(1*R*^{*}, 5*S*^{*})-3-(5-morpholin-4-yl-5-oxopentanoyl)-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid (2-chlorobenzyl)cyclopropylamide;
- 25 (*rac.*)-(1*R*^{*}, 5*S*^{*})-3-(2-tetrazol-1-ylacetyl)-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid (2-chlorobenzyl)cyclopropylamide;
- 30 1:1-mixture of (*1R*, 5*S*)-3-((2*S*)-2-amino-3-hydroxypropionyl)-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid (2-chlorobenzyl)cyclopropylamide and (*1S*, 5*R*)-3-((2*S*)-2-amino-3-

hydroxypropionyl)-7-{4-[3-(2,3,6-trifluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid (2-chlorobenzyl)cyclopropylamide;

5 1:1-mixture of (*1R*, *5S*)-3-((2*S*)-2-aminopropionyl)-7-{4-[3-(2,3,6-trifluoro-
phenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid (2-
chlorobenzyl)cyclopropylamide and (*1S*, *5R*)-3-((2*S*)-2-aminopropionyl)-7-{4-[3-
(2,3,6-trifluoro-phenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-
carboxylic acid (2-chlorobenzyl)cyclopropylamide;

10 (*rac.*)-(1*R**, *5S**)-3-acetyl-7-{4-[2-(2,6-dichloro-4-methylphenoxy)ethoxy]-
phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid cyclopropyl-(2,3-
dichlorobenzyl)amide;

15 (*rac.*)-(1*R**, *5S**)-3-acetyl-7-{4-[2-(2,6-dichloro-4-methylphenoxy)ethoxy]-
phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid cyclopropyl-(2,3-
dimethylbenzyl)amide;

20 (*rac.*)-(1*R**, *5S**)-3-acetyl-7-{4-[2-(2,6-dichloro-4-methylphenoxy)ethoxy]-
phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid (2-chloro-3-
trifluoromethyl-benzyl)cyclopropylamide;

25 (*rac.*)-(1*R**, *5S**)-3-acetyl-7-{4-[2-(2,6-dichloro-4-methylphenoxy)ethoxy]-
phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid (2-bromobenzyl)-
cyclopropylamide;

(*rac.*)-(1*R**, *5S**)-3-acetyl-7-{4-[2-(2,6-dichloro-4-methylphenoxy)ethoxy]-
phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid (2-chlorobenzyl)-
cyclopropylamide;

30 (*rac.*)-(1*R**, *5S**)-3-acetyl-7-{4-[2-(2,6-dichloro-4-methylphenoxy)ethoxy]-
phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid (2-chlorobenzyl)-
ethylamide;

- (*rac.*)-(1*R*^{*}, 5*S*^{*})-3-acetyl-7-{4-[2-(2,6-dichloro-4-methylphenoxy)ethoxy]-phenyl}-3,9-diaza-bicyclo[3.3.1]non-6-ene-6-carboxylic acid (6-chlorobenzo-[1,3]dioxol-5-ylmethyl)cyclopropylamide;
- 5 (*rac.*)-(1*R*^{*}, 5*S*^{*})-3-acetyl-7-{4-[2-(2,6-dichloro-4-methylphenoxy)ethoxy]-phenyl}-3,9-diaza-bicyclo[3.3.1]non-6-ene-6-carboxylic acid cyclopropyl-(3,5-dimethoxybenzyl)amide;
- 10 (*rac.*)-(1*R*^{*}, 5*S*^{*})-3-acetyl-7-{4-[2-(2,6-dichloro-4-methylphenoxy)ethoxy]-phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid cyclopropyl-(3-methoxybenzyl)amide;
- 15 (*rac.*)-(1*R*^{*}, 5*S*^{*})-3-acetyl-7-{4-[2-(2-chloro-5-fluorophenoxy)ethoxy]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid cyclopropyl-(2,3-dichlorobenzyl)amide;
- 20 (*rac.*)-(1*R*^{*}, 5*S*^{*})-3-acetyl-7-{4-[2-(2-chloro-5-fluorophenoxy)ethoxy]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid (2-chloro-3-trifluoromethylbenzyl)cyclopropylamide;
- 25 (*rac.*)-(1*R*^{*}, 5*S*^{*})-3-acetyl-7-{4-[2-(2,6-dichloro-4-methylphenoxy)ethoxy]-phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid (3-chlorobenzyl)cyclopropylamide;
- (*rac.*)-(1*R*^{*}, 5*S*^{*})-3-acetyl-7-{4-[2-(2,6-dichloro-4-methylphenoxy)ethoxy]-phenyl}-3,9-diaza-bicyclo[3.3.1]non-6-ene-6-carboxylic acid cyclopropyl-(3-methylbenzyl)amide;
- 30 (*rac.*)-(1*R*^{*}, 5*S*^{*})-3-acetyl-7-{4-[2-(2,6-dichloro-4-methylphenoxy)ethoxy]-phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid cyclopropyl-(2-fluoro-5-methoxybenzyl)amide;

(*rac.*)-(1*R*^{*}, 5*S*^{*})-3-acetyl-7-{4-[2-(2-chloro-4,5-dimethylphenoxy)ethoxy]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid cyclopropyl-(2,3-dichlorobenzyl)amide;

5

(*rac.*)-(1*R*^{*}, 5*S*^{*})-3-acetyl-7-{4-[3-(2-chloro-3,6-difluorophenoxy)propyl]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid (2-chlorobenzyl)cyclopropylamide;

10 (*rac.*)-(1*R*^{*}, 5*S*^{*})-3-acetyl-7-{4-[2-(2-chloro-4,5-dimethylphenoxy)ethoxy]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid (2-chloro-3-trifluoromethylbenzyl)cyclopropylamide;

15 (*rac.*)-(1*R*^{*}, 5*S*^{*})-3-acetyl-7-{4-[2-(2,6-dichloro-4-methylphenoxy)ethoxy]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid cyclopropyl-(3,4-dimethoxybenzyl)amide;

20 (*rac.*)-(1*R*^{*}, 5*S*^{*})-3-acetyl-7-{4-[2-(2,6-dichloro-4-methylphenoxy)ethoxy]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid cyclopropyl-(3-trifluoromethoxybenzyl)amide;

(*rac.*)-(1*R*^{*}, 5*S*^{*})-3-acetyl-7-{4-[2-(2,4,5-trichlorophenoxy)ethoxy]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid (2-chloro-3-trifluoromethylbenzyl)cyclopropylamide;

25

(*rac.*)-(1*R*^{*}, 5*S*^{*})-3-acetyl-7-{4-[2-(2,4,5-trichlorophenoxy)ethoxy]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid cyclopropyl-(2,3-dichlorobenzyl)amide;

30 (*rac.*)-(1*R*^{*}, 5*S*^{*})-3-acetyl-7-{4-[2-(2,3-dichlorophenoxy)ethoxy]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid cyclopropyl-(2,3-dimethylbenzyl)amide;

(*rac.*)-(1*R*^{*}, 5*S*^{*})-3-acetyl-7-{4-[2-(2,4,5-trichlorophenoxy)ethoxy]phenyl}-3,9-diazabicyclo-[3.3.1]non-6-ene-6-carboxylic acid cyclopropyl-(2,3-dimethylbenzyl)amide;

5

(*rac.*)-(1*R*^{*}, 5*S*^{*})-3-acetyl-7-{4-[2-(2-chloro-4,5-dimethylphenoxy)ethoxy]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid cyclopropyl-(2,3-dimethylbenzyl)amide;

10 (*rac.*)-(1*R*^{*}, 5*S*^{*})-3-acetyl-7-{4-[2-(2-bromo-5-fluorophenoxy)ethoxy]phenyl}-3,9-diaza-bicyclo[3.3.1]non-6-ene-6-carboxylic acid cyclopropyl-(2,3-dichlorobenzyl)amide;

15 (*rac.*)-(1*R*^{*}, 5*S*^{*})-3-acetyl-7-{4-[2-(2-chloro-4,5-dimethylphenoxy)ethoxy]phenyl}-3,9-diaza-bicyclo[3.3.1]non-6-ene-6-carboxylic acid (2-bromobenzyl)cyclopropylamide;

20 (*rac.*)-(1*R*^{*}, 5*S*^{*})-3-acetyl-7-{4-[2-(2,3-dichlorophenoxy)ethoxy]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid cyclopropyl-(2,3-dichlorobenzyl)amide;

(*rac.*)-(1*R*^{*}, 5*S*^{*})-3-acetyl-7-{4-[2-(2-chloro-5-fluorophenoxy)ethoxy]phenyl}-3,9-diazabicyclo[3.3.1]non-6-ene-6-carboxylic acid cyclopropyl-(2,3-dimethylbenzyl)amide;

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(*rac.*)-(1*R*, 5*S*)-7-{4-[3-(2,3,6-trifluorophenoxy)propylphenyl]-3,9-diazabicyclo-[3.3.1]non-6-ene-6-carboxylic acid (2-chlorobenzyl)cyclopropylamide;

30 (*rac.*)-(1*S*, 5*R*)-7-{4-[3-(2,3,6-trifluorophenoxy)propylphenyl]-3,9-diazabicyclo-[3.3.1]non-6-ene-6-carboxylic acid (2-chlorobenzyl)cyclopropylamide.

9. Pharmaceutical compositions containing a compound of any one of claims 1 - 8 and usual carrier materials and adjuvants for the treatment or prophylaxis of disorders which are associated with a dysregulation of the renin-angiotensin system (RAS), comprising cardiovascular and renal diseases hypertension,
5 congestive heart failure, pulmonary hypertension, cardiac insufficiency, renal insufficiency, renal or myocardial ischemia, atherosclerosis, renal failure, erectile dysfunction, glomerulonephritis, renal colic, glaucoma, diabetic complications, complications after vascular or cardiac surgery, restenosis, complications of treatment with immunosuppressive agents after organ transplantation, and other
10 diseases known to be related to the RAS.
 10. A method for the treatment or prophylaxis of diseases which are related to the RAS comprising hypertension, congestive heart failure, pulmonary hypertension, cardiac insufficiency, renal insufficiency, renal or myocardial ischemia,
15 atherosclerosis, renal failure, erectile dysfunction, glomerulonephritis, renal colic, glaucoma, diabetic complications, complications after vascular or cardiac surgery, restenosis, complications of treatment with immunosuppressive agents after organ transplantation, and other diseases which are related to the RAS, which method comprises administrating a compound according to any one of claims 1 to 8 to a
20 human being or animal.
 11. The use of compounds according to any one of claims 1 to 8 for the treatment or prophylaxis of diseases which are associated with the RAS comprising hypertension, congestive heart failure, pulmonary hypertension, cardiac
25 insufficiency, renal insufficiency, renal or myocardial ischemia, atherosclerosis, renal failure, erectile dysfunction, glomerulonephritis, renal colic, glaucoma, diabetic complications, complications after vascular or cardiac surgery, restenosis, complications of treatment with immunosuppressive agents after organ transplantation, and other diseases known to be related to the RAS.
- 30
12. The use of one or more compounds of any one of claims 1 to 8 in combination with other pharmacologically active compounds comprising ACE

inhibitors, angiotensin II receptor antagonists, endothelin receptor antagonists, vasodilators, calcium antagonists, potassium activators, diuretics, sympatholitics, beta-adrenergic antagonists, alpha-adrenergic antagonists, for the treatment of disorders as set forth in any one of claims 9 to 11.

INTERNATIONAL SEARCH REPORT

International Application No PCT/EP 03/03721

A. CLASSIFICATION OF SUBJECT MATTER IPC 7 C07D471/08 A61K31/4995
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According to International Patent Classification (IPC) or to both national classification and IPC

B. FIELDS SEARCHED

Minimum documentation searched (classification system followed by classification symbols) IPC 7 C07D A61K
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Documentation searched other than minimum documentation to the extent that such documents are included in the fields searched

Electronic data base consulted during the international search (name of data base and, where practical, search terms used)
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EPO-Internal, WPI Data, PAJ, BIOSIS, EMBASE, CHEM ABS Data
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C. DOCUMENTS CONSIDERED TO BE RELEVANT

Category	Citation of document, with indication, where appropriate, of the relevant passages	Relevant to claim No.
A	US 3 509 161 A (DOLD OTTO ET AL) 28 April 1970 (1970-04-28) Formula (I); examples 1-9; Table in col. 8 ---	1-12
A	WO 97 09311 A (HOFFMANN LA ROCHE) 13 March 1997 (1997-03-13) cited in the application Claims; Formula (I); examples 82-85, 86(eee), 147 --- -/-	1-12

<input checked="" type="checkbox"/> Further documents are listed in the continuation of box C.
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<input checked="" type="checkbox"/> Patent family members are listed in annex.
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* Special categories of cited documents :

- *A* document defining the general state of the art which is not considered to be of particular relevance
- *E* earlier document but published on or after the international filing date
- *L* document which may throw doubts on priority claim(s) or which is cited to establish the publication date of another citation or other special reason (as specified)
- *O* document referring to an oral disclosure, use, exhibition or other means
- *P* document published prior to the international filing date but later than the priority date claimed

T later document published after the international filing date or priority date and not in conflict with the application but cited to understand the principle or theory underlying the invention

X document of particular relevance; the claimed invention cannot be considered novel or cannot be considered to involve an inventive step when the document is taken alone

Y document of particular relevance; the claimed invention cannot be considered to involve an inventive step when the document is combined with one or more other such documents, such combination being obvious to a person skilled in the art

& document member of the same patent family

Date of the actual completion of the international search

Date of mailing of the international search report
--

4 August 2003

20/08/2003

Name and mailing address of the ISA

European Patent Office, P.B. 5818 Patentlaan 2 NL - 2280 HV Rijswijk Tel. (+31-70) 340-2040, Tx. 31 651 epo nl, Fax (+31-70) 340-3016
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Authorized officer

Rivat, C

INTERNATIONAL SEARCH REPORT

International Application No
PCT/EP 03/03721

C.(Continuation) DOCUMENTS CONSIDERED TO BE RELEVANT

Category	Citation of document, with indication, where appropriate, of the relevant passages	Relevant to claim No.
A	S.A. EL-ABBADY: "ENHANCED REACTIVITY OF PYRIDIN-3-OL TOWARDS 4-PHENYL-1,2,4-TRIAZOLINE-3,5-DIONE: FMO TREATMENT OF THE CYCLOADDITION PROCESS BY ASED-MO CALCULATIONS METHOD" ZAGAZIG J. PHARM. SCI., vol. 5, no. 1, 1996, pages 68-74, XP001154254 Compounds (4) and (5)	1-12

INTERNATIONAL SEARCH REPORTInternational application No.
PCT/EP 03/03721**Box I Observations where certain claims were found unsearchable (Continuation of item 1 of first sheet)**

This International Search Report has not been established in respect of certain claims under Article 17(2)(a) for the following reasons:

1. Claims Nos.: because they relate to subject matter not required to be searched by this Authority, namely:

Although claims 10–12 are directed to a method of treatment of the human/animal body, the search has been carried out and based on the alleged effects of the compound/composition.
2. Claims Nos.: because they relate to parts of the International Application that do not comply with the prescribed requirements to such an extent that no meaningful International Search can be carried out, specifically:
3. Claims Nos.: because they are dependent claims and are not drafted in accordance with the second and third sentences of Rule 6.4(a).

Box II Observations where unity of invention is lacking (Continuation of item 2 of first sheet)

This International Searching Authority found multiple inventions in this international application, as follows:

1. As all required additional search fees were timely paid by the applicant, this International Search Report covers all searchable claims.
2. As all searchable claims could be searched without effort justifying an additional fee, this Authority did not invite payment of any additional fee.
3. As only some of the required additional search fees were timely paid by the applicant, this International Search Report covers only those claims for which fees were paid, specifically claims Nos.:
4. No required additional search fees were timely paid by the applicant. Consequently, this International Search Report is restricted to the invention first mentioned in the claims; it is covered by claims Nos.:

Remark on Protest

- The additional search fees were accompanied by the applicant's protest.
 No protest accompanied the payment of additional search fees.

INTERNATIONAL SEARCH REPORT

Information on patent family members

International Application No
PCT/EP 03/03721

Patent document cited in search report	Publication date	Patent family member(s)	Publication date
US 3509161	A 28-04-1970	NONE	
WO 9709311	A 13-03-1997	AT 242213 T 15-06-2003 AU 708616 B2 05-08-1999 .AU 6743296 A 27-03-1997 BR 9610385 A 06-07-1999 CA 2230931 A1 13-03-1997 CN 1202152 A 16-12-1998 CZ 9800684 A3 14-10-1998 DE 59610509 D1 10-07-2003 WO 9709311 A1 13-03-1997 EP 0863875 A1 16-09-1998 HU 9900926 A2 28-09-1999 JP 11500447 T 12-01-1999 NO 980954 A 28-04-1998 NZ 315677 A 28-02-2000 PL 325425 A1 20-07-1998 RU 2167865 C2 27-05-2001 TR 9800409 T1 21-05-1998 TW 474932 B 01-02-2002 US 6051712 A 18-04-2000 US 6150526 A 21-11-2000 ZA 9607424 A 07-03-1997	